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* * * * * Welcome to STN International * * * * *

NEWS	1		Web Page for STN Seminar Schedule - N. America
NEWS	2	APR 04	STN AnaVist, Version 1, to be discontinued
NEWS	3	APR 15	WPIDS, WPINDEX, and WPIX enhanced with new predefined hit display formats
NEWS	4	APR 28	EMBASE Controlled Term thesaurus enhanced
NEWS	5	APR 28	IMSRESEARCH reloaded with enhancements
NEWS	6	MAY 30	INPAFAMDB now available on STN for patent family searching
NEWS	7	MAY 30	DGENE, PCTGEN, and USGENE enhanced with new homology sequence search option
NEWS	8	JUN 06	EPFULL enhanced with 260,000 English abstracts
NEWS	9	JUN 06	KOREAPAT updated with 41,000 documents
NEWS	10	JUN 13	USPATFULL and USPAT2 updated with 11-character patent numbers for U.S. applications
NEWS	11	JUN 19	CAS REGISTRY includes selected substances from web-based collections
NEWS	12	JUN 25	CA/CAPLUS and USPAT databases updated with IPC reclassification data
NEWS	13	JUN 30	AEROSPACE enhanced with more than 1 million U.S. patent records
NEWS	14	JUN 30	EMBASE, EMBAL, and LEMBASE updated with additional options to display authors and affiliated organizations
NEWS	15	JUN 30	STN on the Web enhanced with new STN AnaVist Assistant and BLAST plug-in
NEWS	16	JUN 30	STN AnaVist enhanced with database content from EPFULL
NEWS	17	JUL 28	CA/CAPLUS patent coverage enhanced
NEWS	18	JUL 28	EPFULL enhanced with additional legal status information from the EPOLINE Register
NEWS	19	JUL 28	IFICDB, IFIPAT, and IFIUDB reloaded with enhancements
NEWS	20	JUL 28	STN Viewer performance improved
NEWS	21	AUG 01	INPADOCDB and INPAFAMDB coverage enhanced
NEWS	22	AUG 13	CA/CAPLUS enhanced with printed Chemical Abstracts page images from 1967-1998
NEWS	23	AUG 15	CAOLD to be discontinued on December 31, 2008
NEWS	24	AUG 15	CAPLUS currency for Korean patents enhanced
NEWS	25	AUG 25	CA/CAPLUS, CASREACT, and IFI and USPAT databases enhanced for more flexible patent number searching
NEWS	26	AUG 27	CAS definition of basic patents expanded to ensure comprehensive access to substance and sequence information

NEWS EXPRESS JUNE 27 08 CURRENT WINDOWS VERSION IS V8.3,
AND CURRENT DISCOVER FILE IS DATED 23 JUNE 2008.

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NEWS IPC8 For general information regarding STN implementation of IPC 8

Enter NEWS followed by the item number or name to see news on that specific topic.

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* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 08:52:59 ON 08 SEP 2008

=> file reg		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	0.21	0.21

FILE 'REGISTRY' ENTERED AT 08:53:08 ON 08 SEP 2008
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STRUCTURE FILE UPDATES: 7 SEP 2008 HIGHEST RN 1047406-12-1
DICTIONARY FILE UPDATES: 7 SEP 2008 HIGHEST RN 1047406-12-1

New CAS Information Use Policies, enter HELP USAGETERMS for details.

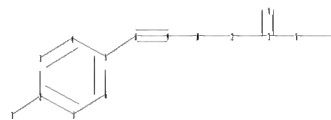
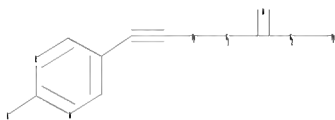
TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

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REGISTRY includes numerically searchable data for experimental and predicted properties as well as tags indicating availability of experimental property data in the original document. For information on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

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Uploading C:\Program Files\STNEXP\Queries\10596745s3.str



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chain nodes :
7 8 9 10 12 13 14 16 17
ring nodes :
1 2 3 4 5 6
chain bonds :
2-7 5-8 8-9 9-10 10-12 12-13 13-14 13-16 16-17
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6
exact/norm bonds :
2-7 9-10 10-12 12-13 13-14 13-16 16-17
exact bonds :
5-8 8-9
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6

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G1:C,O,N

G2:C,O,N

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Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:Atom
12:CLASS 13:CLASS 14:CLASS 16:CLASS 17:Atom

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L1 STRUCTURE UPLOADED

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=> s l1 sss sam
SAMPLE SEARCH INITIATED 08:53:35 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 108 TO ITERATE

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100.0% PROCESSED          108 ITERATIONS                      9 ANSWERS
SEARCH TIME: 00.00.01

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FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**

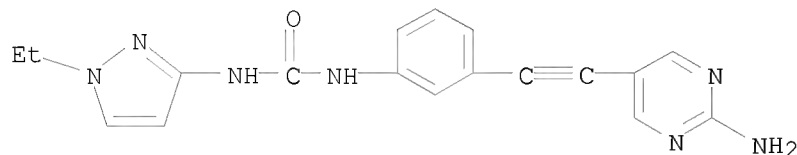
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PROJECTED ITERATIONS: 1537 TO 2783
PROJECTED ANSWERS: 9 TO 360

L2 9 SEA SSS SAM L1

=> d scan

L2 9 ANSWERS REGISTRY COPYRIGHT 2008 ACS on STN
IN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(1-ethyl-1H-
pyrazol-3-yl)-
MF C18 H17 N7 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):end

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	0.92	1.13

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FILE COVERS 1907 - 8 Sep 2008 VOL 149 ISS 11
FILE LAST UPDATED: 7 Sep 2008 (20080907/ED)

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Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

<http://www.cas.org/legal/infopolicy.html>

=> s 12

L3 1 L2

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                               ENTRY      SESSION
FULL ESTIMATED COST          0.48      1.61
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STRUCTURE FILE UPDATES:    7 SEP 2008  HIGHEST RN 1047406-12-1
DICTIONARY FILE UPDATES:   7 SEP 2008  HIGHEST RN 1047406-12-1
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New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH July 5, 2008.

Please note that search-term pricing does apply when
conducting SmartSELECT searches.

REGISTRY includes numerically searchable data for experimental and
predicted properties as well as tags indicating availability of
experimental property data in the original document. For information
on property searching in REGISTRY, refer to:

<http://www.cas.org/support/stngen/stndoc/properties.html>

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FULL SEARCH INITIATED 08:54:43 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -      1944 TO ITERATE
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100.0% PROCESSED      1944 ITERATIONS      172 ANSWERS
SEARCH TIME: 00.00.01
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L4      172 SEA SSS FUL L1
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                               ENTRY      SESSION
FULL ESTIMATED COST          178.36     179.97
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FILE 'CAPLUS' ENTERED AT 08:54:47 ON 08 SEP 2008
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FILE COVERS 1907 - 8 Sep 2008  VOL 149 ISS 11
FILE LAST UPDATED: 7 Sep 2008  (20080907/ED)
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Caplus now includes complete International Patent Classification (IPC) reclassification data for the second quarter of 2008.

Effective October 17, 2005, revised CAS Information Use Policies apply. They are available for your review at:

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=> s 14

L5 3 L4

=> d ibib abs hitstr 1-3

L5 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2006:790854 CAPLUS

DOCUMENT NUMBER: 145:230644

TITLE: Preparation of pyrimidine derivatives and their use as Tie2 receptor tyrosine kinase inhibitors

INVENTOR(S): Jones, Clifford David; Luke, Richard William Arthur; Mccoull, William

PATENT ASSIGNEE(S): Astrazeneca AB, Swed.; Astrazeneca Uk Limited

SOURCE: PCT Int. Appl., 168pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

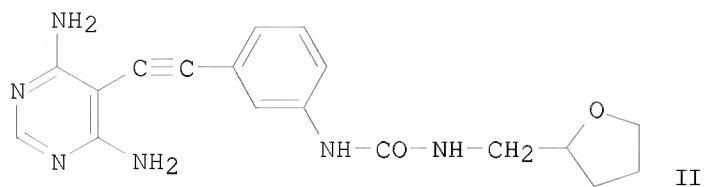
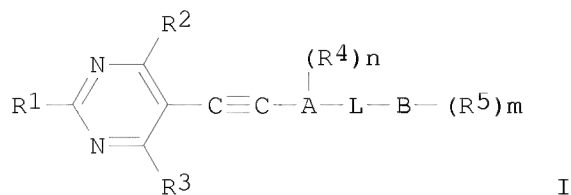
FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

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WO 2006082373	A1	20060810	WO 2006-GB284	20060127
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KM, KN, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, LY, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NG, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SM, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, LV, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
EP 1863805	A1	20071212	EP 2006-701245	20060127
R: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LI, LT, LU, LV, MC, NL, PL, PT, RO, SE, SI, SK, TR				
JP 2008528663	T	20080731	JP 2007-553683	20060127
IN 2007DN05604	A	20070817	IN 2007-DN5604	20070719
CN 101137652	A	20080305	CN 2006-80007855	20070911
PRIORITY APPLN. INFO.:			GB 2005-1984	A 20050201
			GB 2005-2417	A 20050205
			GB 2005-12614	A 20050621
			WO 2006-GB284	W 20060127

OTHER SOURCE(S): MARPAT 145:230644

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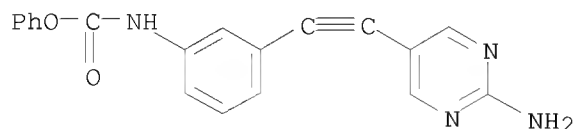


AB Substituted pyrimidine derivs. I, wherein R1 is an (un)substituted amine, (un)substituted 3-7 membered heterocyclic ring; R2 and R3 are H, (un)substituted alkyl, (un)substituted alkoxy; A is a 5 or 6 membered heteroaryl ring; R4 is halo, cyano, alkoxy, cyclopropyl, alkyl, where the alkoxy or alkyl groups are optionally substituted by cyano or 1 or more fluoro groups; L is meta or para attached by an (un)substituted amide, (un)substituted amine, alkyl group; B is a cycloalkyl, heterocyclic ring, aryl, heteroaryl, bicyclic ring; R5 is a halo, hydroxyl, amino, alkylamino, cyano, cycloalkyl ring, an (un)substituted 3 to 7 membered heterocyclic ring; m and n are 0-3 are prepared and used as medicaments and in the production of an anti-angiogenic effect in a warm blooded animal. Thus, II was prepared and tested as an in vitro inhibitor of the Tie2 receptor tyrosine kinase and in the inhibition of autophosphorylation of Tie2 receptor tyrosine kinase (IC₅₀ are 1.5 and 1.9 μ M resp.). Further, I can be used in the treatment of cancer and as antineoplastic prodrugs.

IT 857265-17-9P, Phenyl[3-[(2-aminopyrimidin-5-yl)ethynyl]phenyl]carbamate 857266-46-7P, Phenyl 3-[[2-[[3-(dimethylamino)propyl]amino]pyrimidin-5-yl]ethynyl]phenylcarbamate 905439-39-6P 905439-44-3P 905439-48-7P 905439-61-4P 905439-64-7P
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
 (preparation of pyrimidine derivs. and their use as Tie2 receptor tyrosine kinase inhibitors)

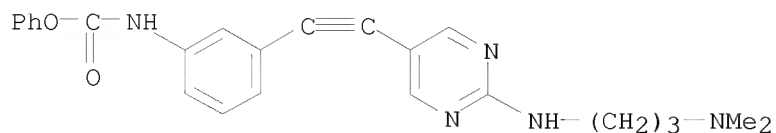
RN 857265-17-9 CAPLUS

CN Carbamic acid, [3-[(2-amino-5-pyrimidinyl)ethynyl]phenyl]-, phenyl ester (9CI) (CA INDEX NAME)



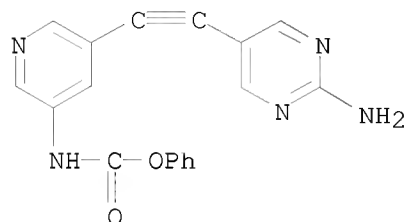
RN 857266-46-7 CAPLUS

CN Carbamic acid, [3-[[2-[[3-(dimethylamino)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-, phenyl ester (9CI) (CA INDEX NAME)



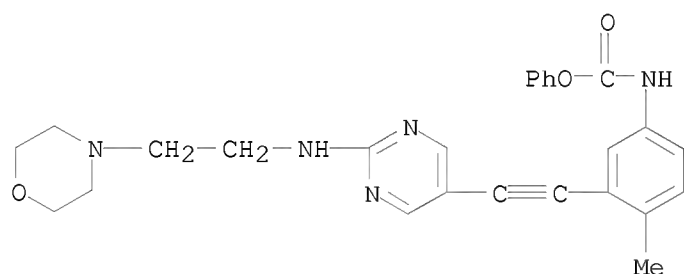
RN 905439-39-6 CAPLUS

CN Carbamic acid, [5-[(2-amino-5-pyrimidinyl)ethynyl]-3-pyridinyl]-, phenyl ester (9CI) (CA INDEX NAME)



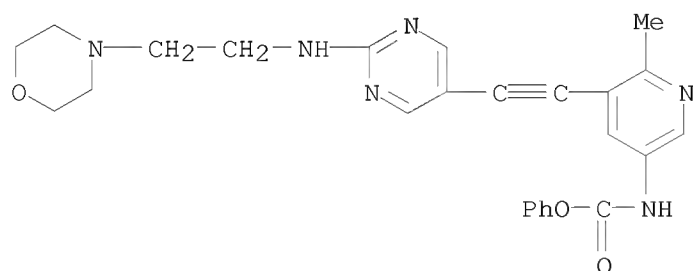
RN 905439-44-3 CAPLUS

CN Carbamic acid, [4-methyl-3-[[2-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-, phenyl ester (9CI) (CA INDEX NAME)



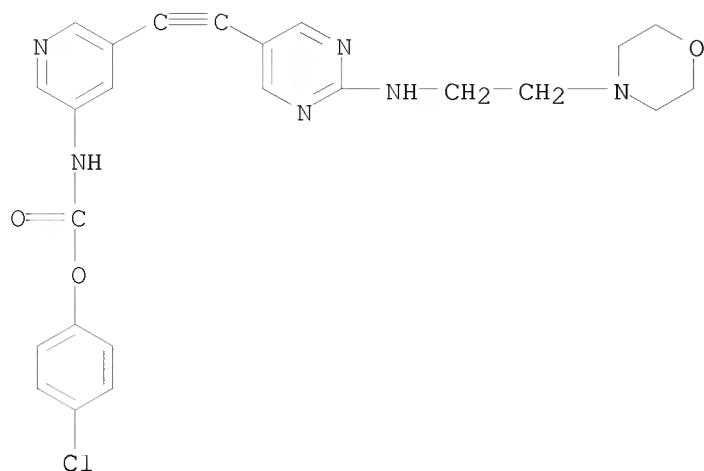
RN 905439-48-7 CAPLUS

CN Carbamic acid, [6-methyl-5-[[2-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]-3-pyridinyl]-, phenyl ester (9CI) (CA INDEX NAME)

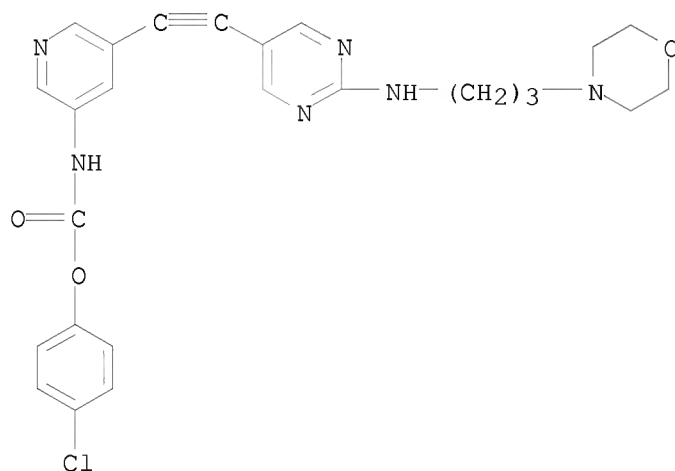


RN 905439-61-4 CAPLUS

CN Carbamic acid, [5-[[2-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]-3-pyridinyl]-, 4-chlorophenyl ester (9CI) (CA INDEX NAME)



RN 905439-64-7 CAPLUS
 CN Carbamic acid, [5-[[2-[[3-(4-morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]-3-pyridinyl]-, 4-chlorophenyl ester (9CI) (CA INDEX NAME)



REFERENCE COUNT: 5 THERE ARE 5 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2008 ACS on STN

ACCESSION NUMBER: 2006:381241 CAPLUS

DOCUMENT NUMBER: 144:432828

TITLE: Heteroaryl-substituted alkyne compounds as protein kinase inhibitors, their preparation, pharmaceutical compositions, and use in therapy

INVENTOR(S): Chaffee, Stuart C.; Albrecht, Brian K.; Hodous, Brian L.; Martin, Matthew W.; McGowan, David C.; Dimauro, Erin F.; Reddy, Gade; Cee, Victor J.; Olivieri, Philip R.; Reed, Anthony; Romero, Karina

PATENT ASSIGNEE(S): Amgen Inc., USA

SOURCE: PCT Int. Appl., 330 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent

LANGUAGE: English

FAMILY ACC. NUM. COUNT: 1

PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2006044823	A2	20060427	WO 2005-US37299	20051017
WO 2006044823	A3	20060615		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KM, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, LY, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NG, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SM, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
RW: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, LV, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG, BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
US 20060217380	A1	20060928	US 2005-251490	20051014
AU 2005295414	A1	20060427	AU 2005-295414	20051017
CA 2583907	A1	20060427	CA 2005-2583907	20051017
EP 1802586	A2	20070704	EP 2005-812237	20051017
R: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LI, LT, LU, LV, MC, NL, PL, PT, RO, SE, SI, SK, TR, AL, BA, HR, MK, YU				
PRIORITY APPLN. INFO.:			US 2004-620100P	P 20041018
			US 2005-251490	A 20051014
			WO 2005-US37299	W 20051017
OTHER SOURCE(S):			MARPAT 144:432828	
GI				

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

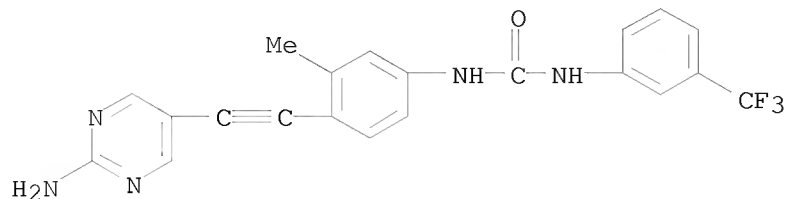
AB The invention relates to heteroaryl-substituted alkynes of formula I, which are protein kinase modulators. In compds. I, W, X, Y, and Z are independently selected from N and (un)substituted C; R1 is (un)substituted amino, acyl, acyloxy, carboxylate, carbamoyl, thiocarbamoyl, etc.; and R2 is 5- to 8-membered monocyclic, 6- to 12-membered bicyclic, or 7- to 14-membered tricyclic ring system, optionally including 1-3 heteroatoms selected from O, N, and S; including stereoisomers, tautomers, solvates, salts, derivs., and prodrugs thereof. The invention also relates to the preparation of I, pharmaceutical compns. comprising a compound I and a pharmaceutically acceptable carrier, as well as to the use of the compns. for the prophylaxis and treatment of protein kinase-mediated diseases, including inflammation, cancer and related conditions. Chlorination of 3-iodo-4-methylbenzoic acid and amidation with 3-trifluoromethylaniline gave benzamide II, which underwent coupling with 2-amino-5-ethynylpyrimidine (preparation from 2-amino-5-iodopyrimidine and trimethylsilylacetylene is given) to give pyrimidinylalkyne III. Several compds. of the invention, e.g., III, express IC50 values of less than or equal to 10 μ M both for Tie-2 and Lck kinase.

IT 884602-98-6P, N-[4-[2-(2-Aminopyrimidin-5-yl)ethynyl]-3-methylphenyl]-N'-[3-(trifluoromethyl)phenyl]urea 884603-00-3P
 884604-13-1P, N-[3-Methyl-4-[[4-[[2-(methyloxy)phenyl]oxy]-2-[[4-(4-methyl-1-piperazinyl)phenyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[3-(trifluoromethyl)phenyl]urea
 RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(drug candidate; preparation of heteroaryl-substituted alkynes as protein kinase modulators)

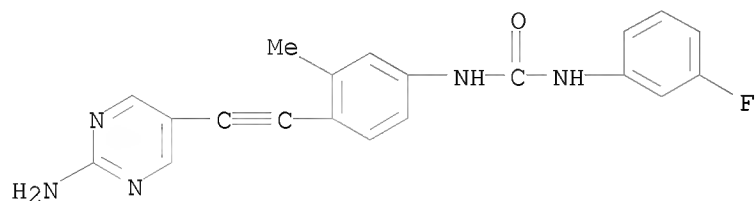
RN 884602-98-6 CAPLUS

CN Urea, N-[4-[2-(2-amino-5-pyrimidinyl)ethynyl]-3-methylphenyl]-N'-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



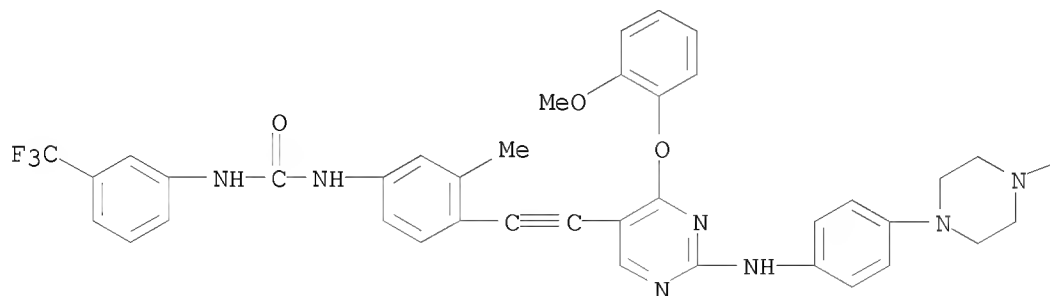
RN 884603-00-3 CAPLUS

CN Urea, N-[4-[2-(2-amino-5-pyrimidinyl)ethynyl]-3-methylphenyl]-N'-(3-fluorophenyl)- (CA INDEX NAME)



RN 884604-13-1 CAPLUS

CN Urea, N-[4-[2-[4-(2-methoxyphenoxy)-2-[[4-(4-methyl-1-piperazinyl)phenyl]amino]-5-pyrimidinyl]ethynyl]-3-methylphenyl]-N'-[3-(trifluoromethyl)phenyl]- (CA INDEX NAME)



PAGE 1-A

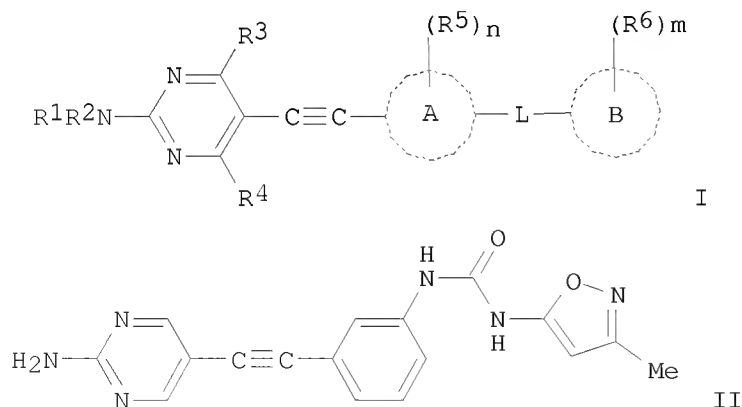
PAGE 1-B

Me

ACCESSION NUMBER: 2005:588668 CAPLUS
 DOCUMENT NUMBER: 143:115557
 TITLE: Preparation of 2-aminopyrimidine derivatives as inhibitors of Tie2 receptor tyrosine kinases
 INVENTOR(S): Jones, Clifford David; Luke, Richard William Arthur; McCoull, William
 PATENT ASSIGNEE(S): Astrazeneca AB, Swed.; Astrazeneca UK Limited
 SOURCE: PCT Int. Appl., 178 pp.
 CODEN: PIXXD2
 DOCUMENT TYPE: Patent
 LANGUAGE: English
 FAMILY ACC. NUM. COUNT: 1
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2005060970	A1	20050707	WO 2004-GB5337	20041220
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
EP 1737463	A1	20070103	EP 2004-806139	20041220
R: AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LI, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR				
CN 1917879	A	20070221	CN 2004-80041901	20041220
JP 2007517007	T	20070628	JP 2006-546306	20041220
US 20080108608	A1	20080508	US 2006-596745	20060622
IN 2006MN00846	A	20070608	IN 2006-MN846	20060717
PRIORITY APPLN. INFO.:				
			GB 2003-30000	A 20031224
			GB 2004-16849	A 20040729
			WO 2004-GB5337	W 20041220

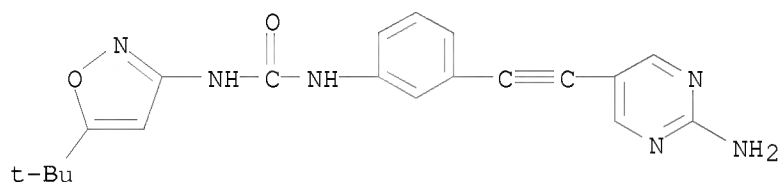
OTHER SOURCE(S): MARPAT 143:115557
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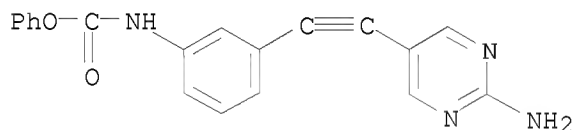
AB Title compds. I [wherein R1, R2 = H, alkyl, alkanoyl; R3, R4 = H, alkyl, alkoxy; R5 = cyclopropyl, halo, cyano; m, n = 0-3; R6 = halo, oxo, cyano; etc., or salts thereof] were prepared as inhibitors of Tie2 receptor

tyrosine kinases. Processes for the synthesis of I and some intermediates involved are claimed. For example, 2-amino-5-iodopyrimidine underwent Pd-catalyzed coupling with 3-ethynylaniline in the presence of CuI. The resultant substituted aniline was condensed with a carbamate, which was obtained from Ph chloroformate and 5-amino-3-methylisoxazole, to give urea II. This compound showed inhibition against Tie2 receptor tyrosine kinase in vitro and inhibition of autophosphorylation of Tie2 receptor tyrosine kinase with IC50 values of 19.871 μ M and 0.337 μ M, resp. Therefore, I and their pharmaceutical compns. have potential use in the production of an anti-angiogenic effect in a warm-blooded animal.

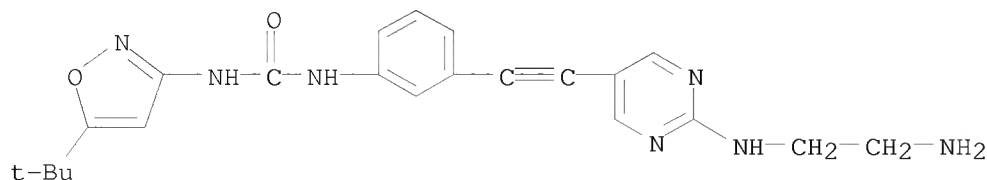
- IT 857265-16-8P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-(5-tert-butylisoxazol-3-yl)urea 857265-17-9P, Phenyl [3-[(2-aminopyrimidin-5-yl)ethynyl]phenyl]carbamate 857265-31-7P, N-[3-[[2-[(2-Aminoethyl)amino]pyrimidin-5-yl]ethynyl]phenyl]-N'-(5-tert-butylisoxazol-3-yl)urea 857265-32-8P, N-[3-[[2-[(3-Aminopropyl)amino]pyrimidin-5-yl]ethynyl]phenyl]-N'-(5-tert-butylisoxazol-3-yl)urea 857266-46-7P, Phenyl [3-[[2-[[3-(dimethylamino)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]carbamate
 RL: PAC (Pharmacological activity); RCT (Reactant); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
 (inhibitor; preparation of pyrimidine derivs. as inhibitors of Tie2 receptor tyrosine kinases)
 RN 857265-16-8 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



- RN 857265-17-9 CAPLUS
 CN Carbamic acid, [3-[(2-amino-5-pyrimidinyl)ethynyl]phenyl]-, phenyl ester (9CI) (CA INDEX NAME)

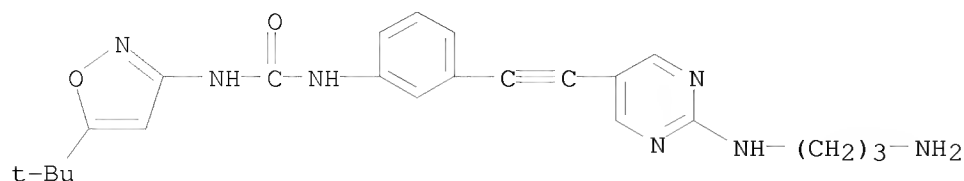


- RN 857265-31-7 CAPLUS
 CN Urea, N-[3-[2-[2-[(2-aminoethyl)amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



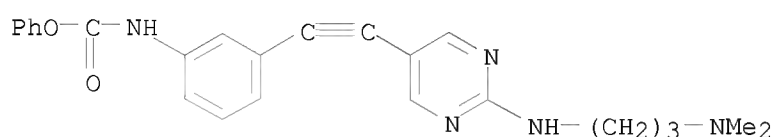
- RN 857265-32-8 CAPLUS

CN Urea, N-[3-[2-[2-[(3-aminopropyl)amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



RN 857266-46-7 CAPLUS

CN Carbamic acid, [3-[[2-[[3-(dimethylamino)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-, phenyl ester (9CI) (CA INDEX NAME)



IT 857264-91-6P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[2-fluoro-5-(trifluoromethyl)phenyl]urea 857264-93-8P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[2-(trifluoromethyl)phenyl]urea 857264-94-9P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[4-(trifluoromethyl)phenyl]urea 857264-95-0P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[2-fluorophenyl]urea 857264-96-1P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[3-fluorophenyl]urea 857264-97-2P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[4-fluorophenyl]urea 857264-98-3P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[3-methoxyphenyl]urea 857264-99-4P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[2,5-difluorophenyl]urea 857265-00-0P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[1,3-benzodioxol-5-yl]urea 857265-01-1P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[3-(trifluoromethyl)phenyl]urea 857265-02-2P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[2-methoxyphenyl]urea 857265-03-3P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[4-methoxyphenyl]urea 857265-04-4P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[3,4-difluorophenyl]urea 857265-05-5P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[3-cyanophenyl]urea 857265-06-6P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[3-chlorophenyl]urea 857265-07-7P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-cyclopentylurea 857265-08-8P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[3,5-difluorophenyl]urea 857265-09-9P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[5-tert-butyl-1,3,4-thiadiazol-2-yl]urea 857265-13-5P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[3-methylisoxazol-5-yl]urea 857265-14-6P, N-[3-[[[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]amino]carbonyl]amino]phenyl]acetamide 857265-15-7P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[4-(trifluoromethyl)pyridin-2-yl]urea 857265-18-0P, N-[3-[(2-Aminopyrimidin-5-yl)ethynyl]phenyl]-N'-[2-oxopiperidin-3-yl]urea 857265-19-1P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-(methylamino)pyrimidin-5-yl]ethynyl]phenyl]urea 857265-22-6P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-(dimethylamino)pyrimidin-5-yl]ethynyl]phenyl]urea 857265-23-7P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-(morpholin-4-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-24-8P, N-(5-tert-Butylisoxazol-3-

yl)-N'-[3-[[2-[[3-(morpholin-4-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-25-9P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[(2-methoxyethyl)amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-26-0P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[3-(1H-imidazol-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-27-1P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[(3-methoxypropyl)amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-28-2P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[(2-hydroxyethyl)amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-29-3P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-30-6P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[3-(pyrrolidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-33-9P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-(dimethylamino)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-34-0P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[3-(dimethylamino)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-35-1P, N-[5-[[3-[[[(5-tert-Butylisoxazol-3-yl)amino]carbonyl]amino]phenyl]ethynyl]pyrimidin-2-yl]glycinamide 857265-36-2P 857265-37-3P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-(1H-imidazol-4-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-38-4P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-(pyridin-2-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-39-5P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[3-(isopropylamino)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-40-8P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[3-(4-methylpiperazin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-41-9P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-(pyridin-4-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-42-0P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-43-1P, N-(5-Methylisoxazol-3-yl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-47-5P, N-(5-tert-Butyl-1,3,4-thiadiazol-2-yl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-48-6P, N-(3-Methylisothiazol-5-yl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-49-7P, N-(3-Fluorophenyl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-50-0P, N-(4-Methoxyphenyl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-51-1P, N-(2-Fluorophenyl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-52-2P, N-(2,5-Difluorophenyl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-53-3P, N-(3,4-Difluorophenyl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-54-4P, N-[2-Fluoro-5-(trifluoromethyl)phenyl]-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-55-5P, N-[3-[[2-[[2-(Pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]-N'-[4-(trifluoromethyl)phenyl]urea 857265-56-6P, N-(1,3-Benzodioxol-5-yl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-57-7P, N-(4-Fluorophenyl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-58-8P, N-(3-Chlorophenyl)-N'-[3-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-59-9P, N-(5-Methylisoxazol-3-yl)-N'-[3-[[2-[[2-(morpholin-4-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-60-2P, N-(5-tert-Butyl-1,3,4-thiadiazol-2-yl)-N'-[3-[[2-[[2-(morpholin-4-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-61-3P, N-[2-Fluoro-5-(trifluoromethyl)phenyl]-N'-[3-[[2-[[2-(morpholin-4-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-62-4P, N-(5-Methylisoxazol-3-yl)-N'-[3-[[2-[[3-(morpholin-4-

yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-63-5P,
N-(5-tert-Butyl-1,3,4-thiadiazol-2-yl)-N'-[3-[[2-[[3-(morpholin-4-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-64-6P,
N-[2-Fluoro-5-(trifluoromethyl)phenyl]-N'-[3-[[2-[[3-(morpholin-4-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-65-7P,
N-(5-Methylisoxazol-3-yl)-N'-[4-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-68-0P,
N-(5-tert-Butylisoxazol-3-yl)-N'-[4-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-69-1P,
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N-[2-Fluoro-5-(trifluoromethyl)phenyl]-N'-[4-[[2-[[2-(pyrrolidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857265-71-5P,
N-[5-[[3-[[[(5-tert-Butylisoxazol-3-yl)amino]carbonyl]amino]phenyl]ethynyl]pyrimidin-2-yl]-2-(2-methoxyethoxy)acetamide 857265-72-6P,
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N-[5-[[2-Aminopyrimidin-5-yl]ethynyl]-1,3,4-thiadiazol-2-yl]-N'-[2-fluoro-5-(trifluoromethyl)phenyl]urea 857265-82-8P,
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N-[3-[[2-Aminopyrimidin-5-yl]ethynyl]phenyl]-2-(3-methylisoxazol-5-yl)acetamide 857265-91-9P, N-[4-[[2-Aminopyrimidin-5-yl]ethynyl]phenyl]-2-(2-methoxyphenyl)acetamide 857265-92-0P,
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N-[3-[[2-Aminopyrimidin-5-yl]ethynyl]phenyl]-N'-(1-tert-butyl-3-cyclopropyl-1H-pyrazol-5-yl)urea 857265-99-7P,
N-[3-[[2-Aminopyrimidin-5-yl]ethynyl]phenyl]-N'-(5-methyl-1,3,4-thiadiazol-2-yl)urea 857266-00-3P, N-[3-[[2-Aminopyrimidin-5-yl]ethynyl]phenyl]-N'-(5-ethyl-1,3,4-thiadiazol-2-yl)urea 857266-01-4P, N-[3-[[2-Aminopyrimidin-5-yl]ethynyl]phenyl]-N'-(5-isopropyl-1,3,4-thiadiazol-2-yl)urea 857266-02-5P,
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N-[3-[[2-Aminopyrimidin-5-yl]ethynyl]phenyl]-N'-[5-(trifluoromethyl)-1,3,4-thiadiazol-2-yl]urea 857266-05-8P, N'-[3-[[2-Aminopyrimidin-5-yl]ethynyl]phenyl]-N-methyl-N-[5-(trifluoromethyl)-1,3,4-thiadiazol-2-yl]urea 857266-06-9P, N-[3-[[2-Aminopyrimidin-5-yl]ethynyl]phenyl]-N'-(5-cyclopropyl-1,3,4-thiadiazol-2-yl)urea 857266-07-0P, N-Phenyl-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-09-2P,
N-(5-Methylisoxazol-3-yl)-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-10-5P,
N-[3-[[2-[[3-(Piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]-

N'-[4-(trifluoromethyl)pyridin-2-yl]urea 857266-11-6P,
N-(5-tert-Butyl-1,3,4-thiadiazol-2-yl)-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-12-7P,
N-(3-Methylisoxazol-5-yl)-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-13-8P,
N-(2-Methoxyphenyl)-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-14-9P, N-(3-Fluorophenyl)-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-15-0P, N-[3-[[2-[[4-Aminobutyl]amino]pyrimidin-5-yl]ethynyl]phenyl]-N'-(5-tert-butylisoxazol-3-yl)urea 857266-16-1P
, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-(piperidin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-17-2P,
N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-(isopropylamino)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-18-3P,
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N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-(piperazin-1-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-27-4P,
N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[3-(piperazin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-29-6P,
N-(5-tert-Butylisoxazol-3-yl)-N-methyl-N'-[3-[[2-[[2-(morpholin-4-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-32-1P,
N-(5-tert-Butylisoxazol-3-yl)-N-methyl-N'-[3-[[2-[[3-(morpholin-4-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-33-2P,
N-(5-tert-Butylisoxazol-3-yl)-N-methyl-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-34-3P,
N-(3-tert-Butyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-36-5P,
N-(3-tert-Butyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[[2-[[3-(morpholin-4-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-37-6P,
N-(3-tert-Butyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[[2-[[2-(morpholin-4-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-38-7P,
N-(3-tert-Butyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[[2-[[2-(dimethylamino)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-39-8P, N-(3-tert-Butyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[[2-[[2-(isopropylamino)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-40-1P, N-(3-Cyclopropyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[[2-[[3-(piperidin-1-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-42-3P, N-(3-Cyclopropyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[[2-[[3-(morpholin-4-yl)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-43-4P, N-(3-Cyclopropyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[[2-[[2-(morpholin-4-yl)ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-44-5P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[2-[[2-(2-hydroxy-1-oxoethyl)amino]ethyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-45-6P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[3-[[2-(2-hydroxyethyl)amino]propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-48-9P, N-[3-[[2-[[3-(Dimethylamino)propyl]amino]pyrimidin-5-yl]ethynyl]phenyl]-N'-phenylurea 857266-49-0P,
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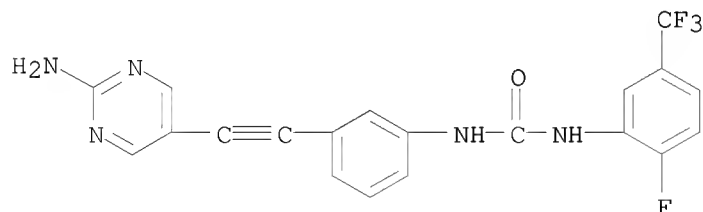
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RL: PAC (Pharmacological activity); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)

(inhibitor; preparation of pyrimidine derivs. as inhibitors of Tie2 receptor tyrosine kinases)

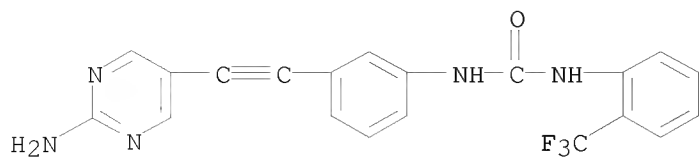
RN 857264-91-6 CAPLUS

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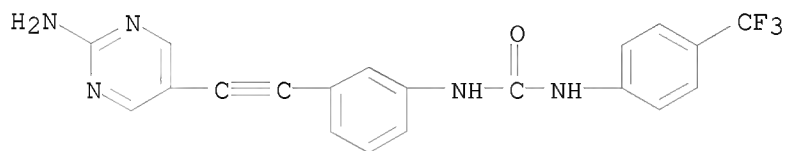
RN 857264-93-8 CAPLUS

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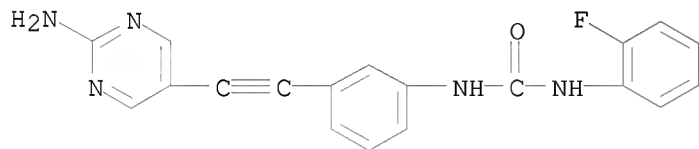
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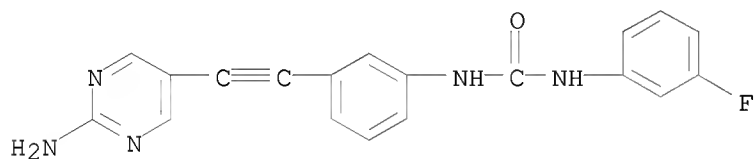
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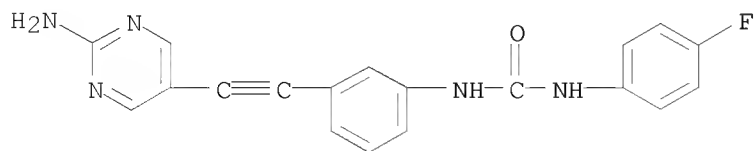
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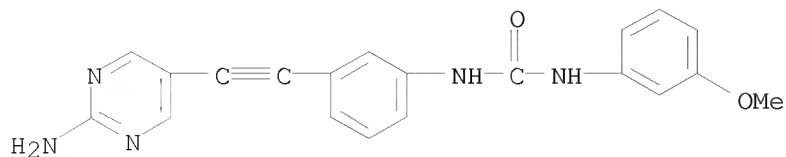
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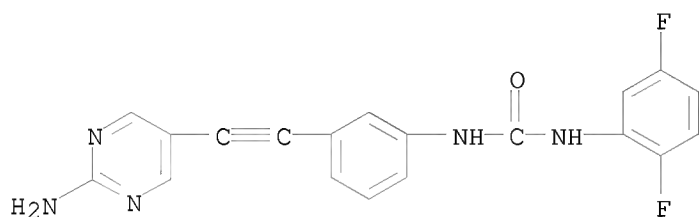


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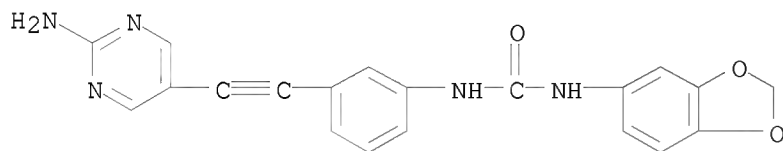
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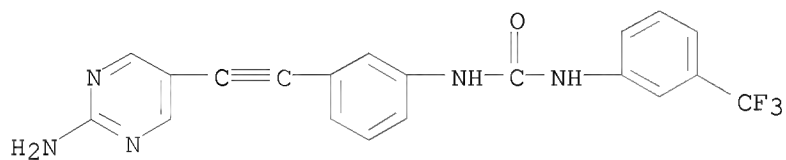
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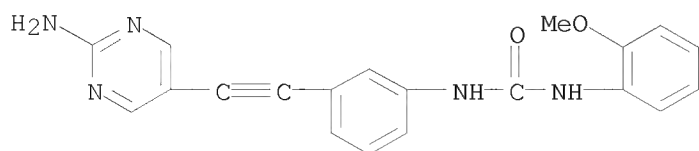
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RN 857265-01-1 CAPLUS
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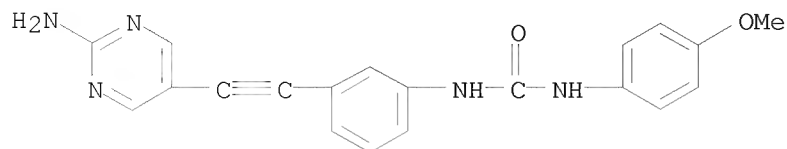


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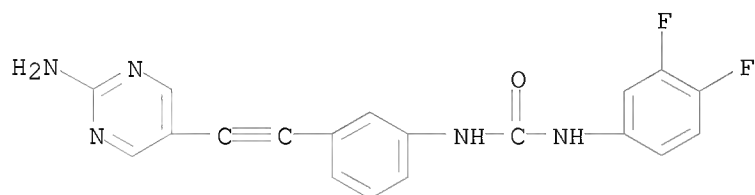
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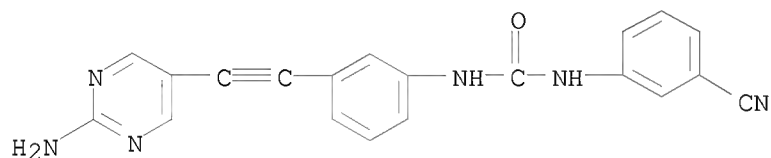
RN 857265-04-4 CAPLUS

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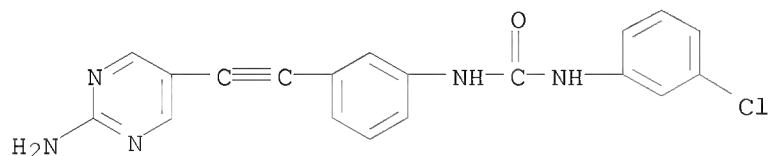
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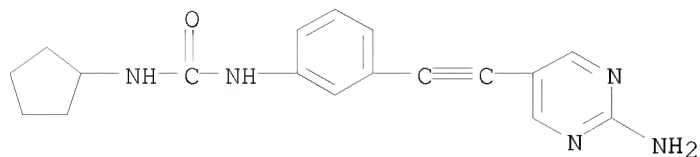
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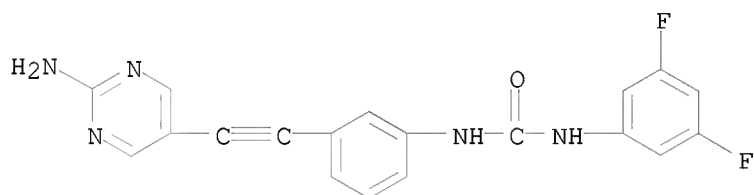


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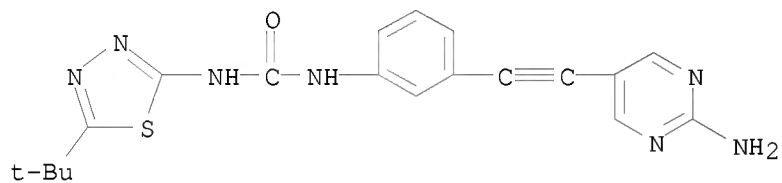
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(CA INDEX NAME)



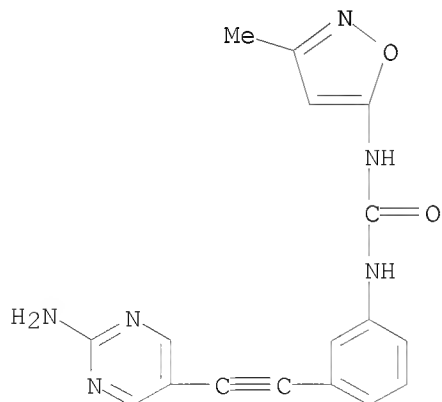
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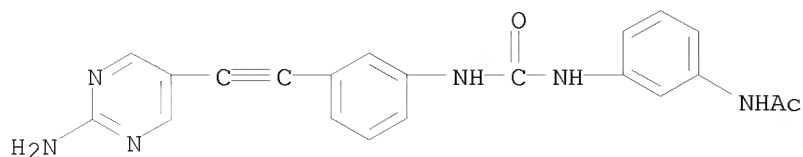
RN 857265-09-9 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl]- (CA INDEX NAME)



RN 857265-13-5 CAPLUS
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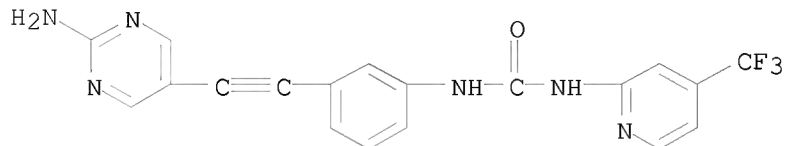


RN 857265-14-6 CAPLUS
 CN Acetamide, N-[3-[[[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]amino]carbonyl]amino]phenyl]- (CA INDEX NAME)



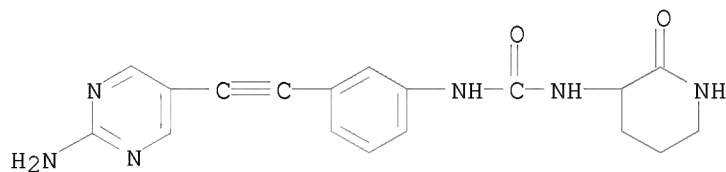
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CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[4-(trifluoromethyl)-2-pyridinyl]- (CA INDEX NAME)



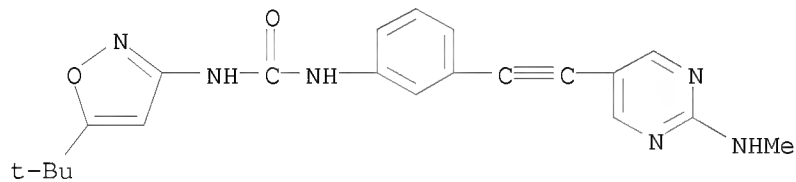
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CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(2-oxo-3-piperidinyl)- (CA INDEX NAME)



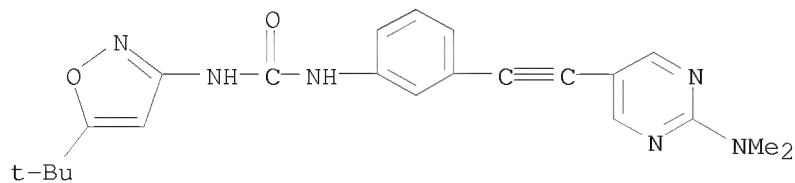
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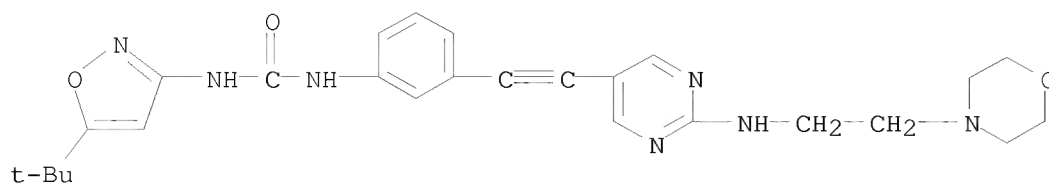
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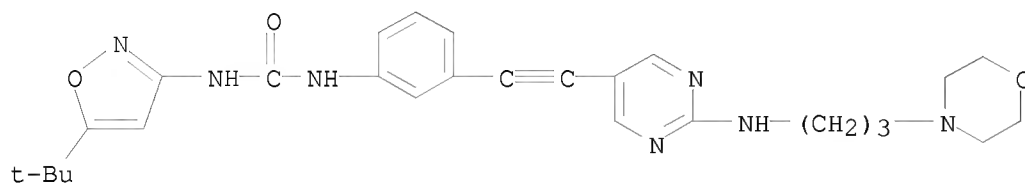
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CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



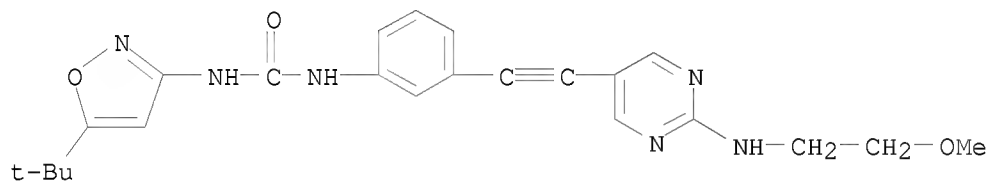
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CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[3-(4-morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



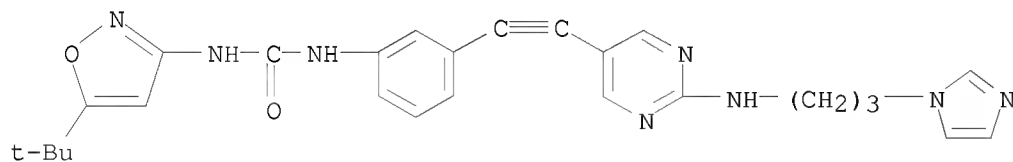
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CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[(2-methoxyethyl)amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



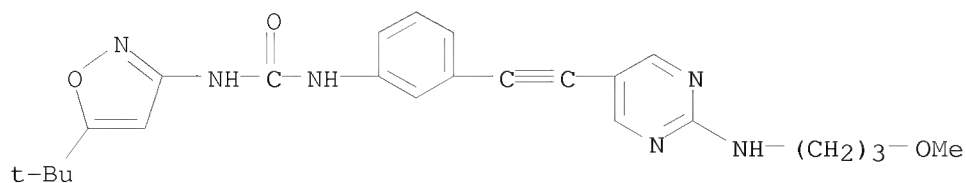
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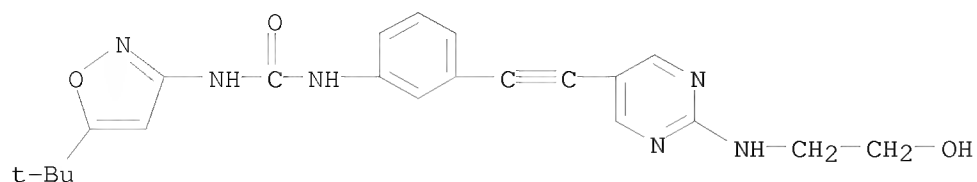
RN 857265-27-1 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[(3-methoxypropyl)amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



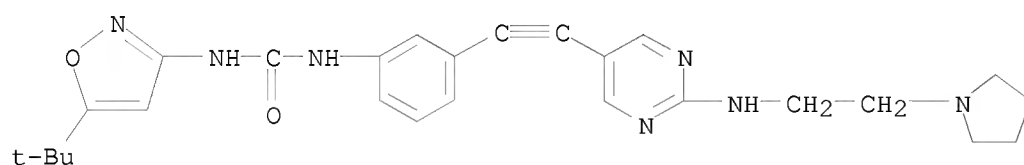
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CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[(2-hydroxyethyl)amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



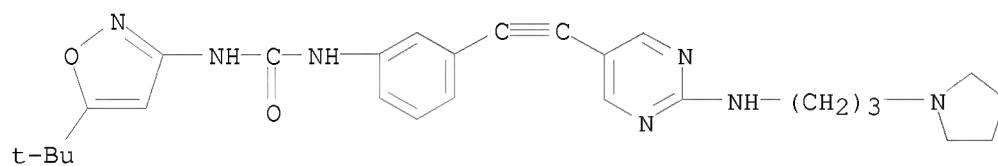
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CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



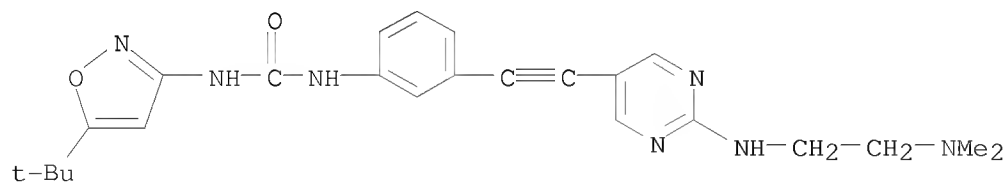
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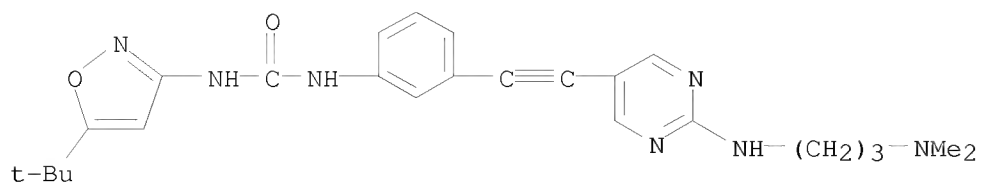
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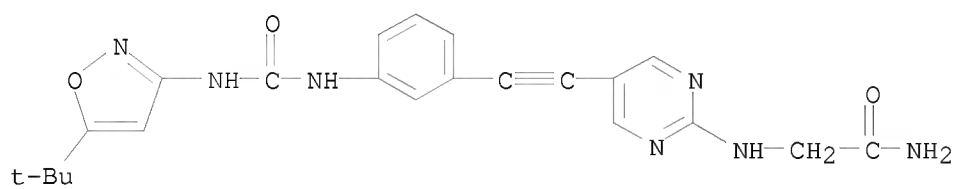
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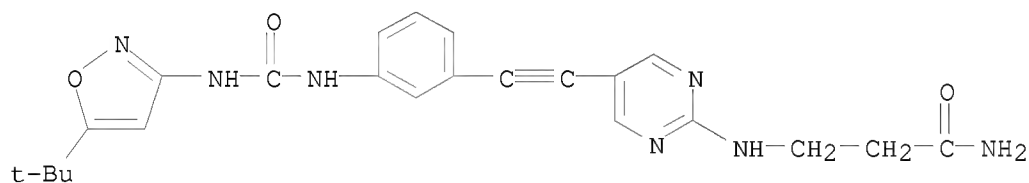
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CN Acetamide, 2-[[5-[2-[3-[[[5-(1,1-dimethylethyl)-3-isoxazolyl]amino]carbonyl]amino]phenyl]ethynyl]-2-pyrimidinyl]amino]- (CA INDEX NAME)



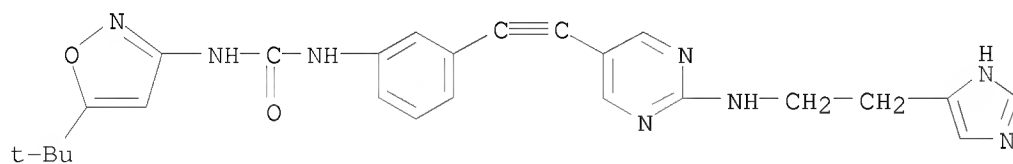
RN 857265-36-2 CAPLUS

CN Propanamide, 3-[[5-[2-[3-[[[5-(1,1-dimethylethyl)-3-isoxazolyl]amino]carbonyl]amino]phenyl]ethynyl]-2-pyrimidinyl]amino]- (CA INDEX NAME)



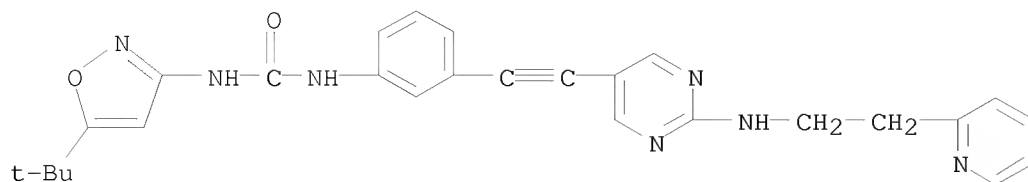
RN 857265-37-3 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[2-(1H-imidazol-5-yl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



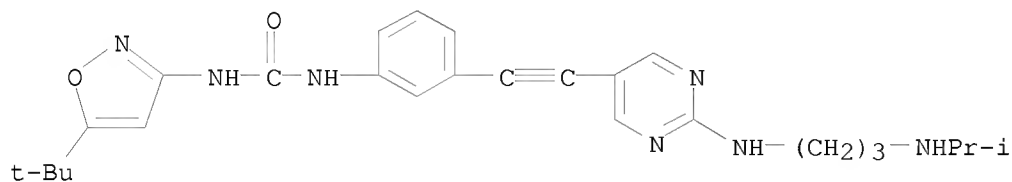
RN 857265-38-4 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[2-(2-pyridinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857265-39-5 CAPLUS

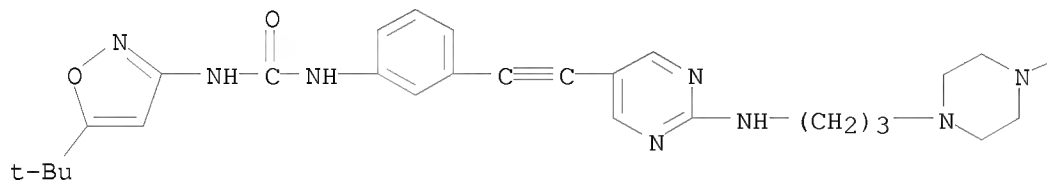
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[3-[(1-methylethyl)amino]propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857265-40-8 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[3-(4-methyl-1-piperazinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

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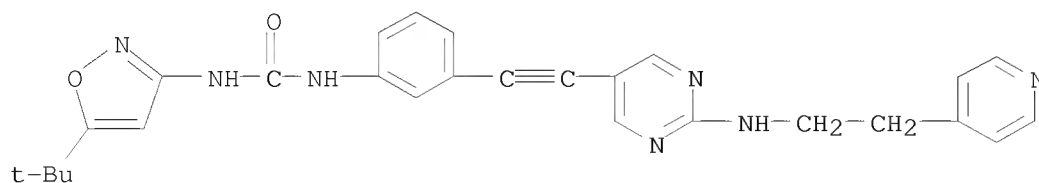


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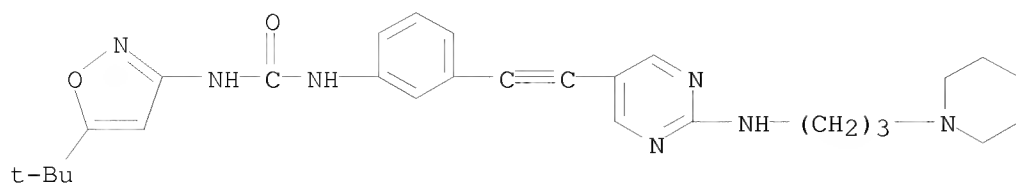
RN 857265-41-9 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[2-(4-pyridinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



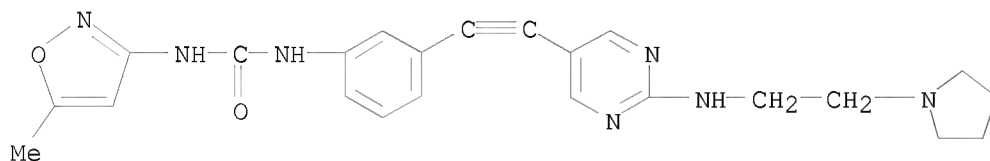
RN 857265-42-0 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[3-(1-piperidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



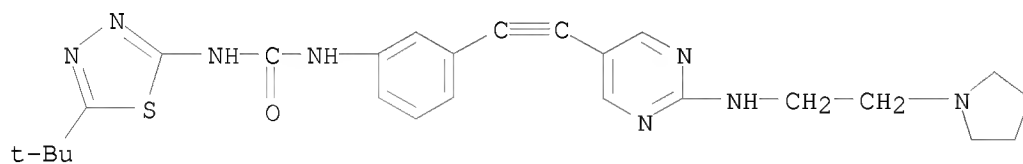
RN 857265-43-1 CAPLUS

CN Urea, N-(5-methyl-3-isoxazolyl)-N'-[3-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



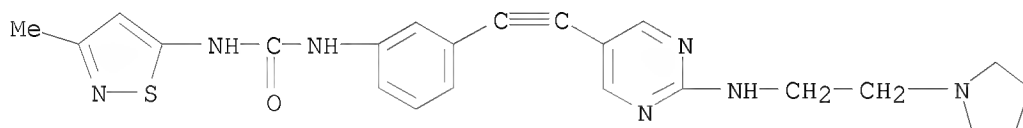
RN 857265-47-5 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl]-N'-[3-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



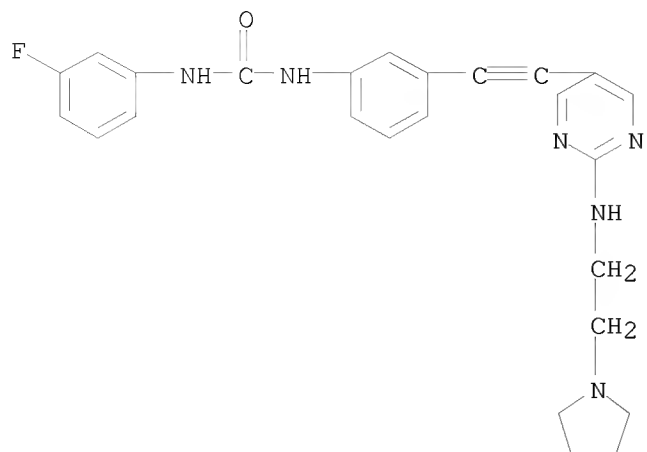
RN 857265-48-6 CAPLUS

CN Urea, N-(3-methyl-5-isothiazolyl)-N'-[3-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



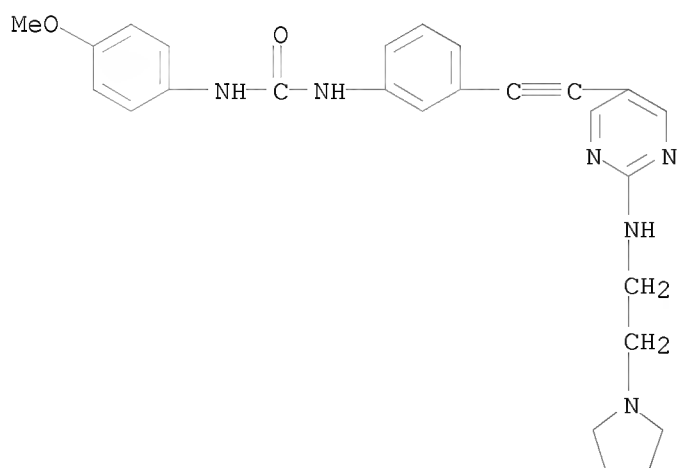
RN 857265-49-7 CAPLUS

CN Urea, N-(3-fluorophenyl)-N'-[3-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



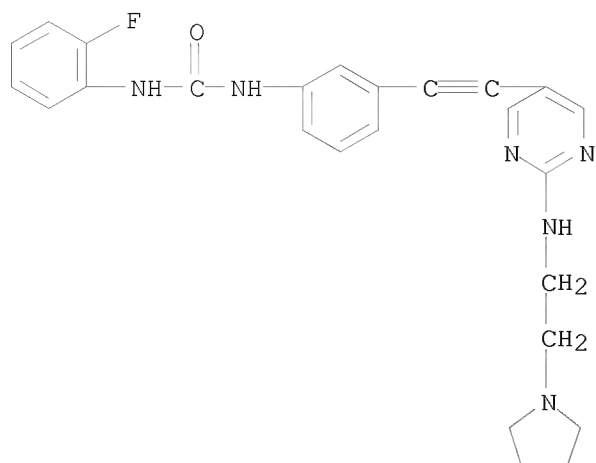
RN 857265-50-0 CAPLUS

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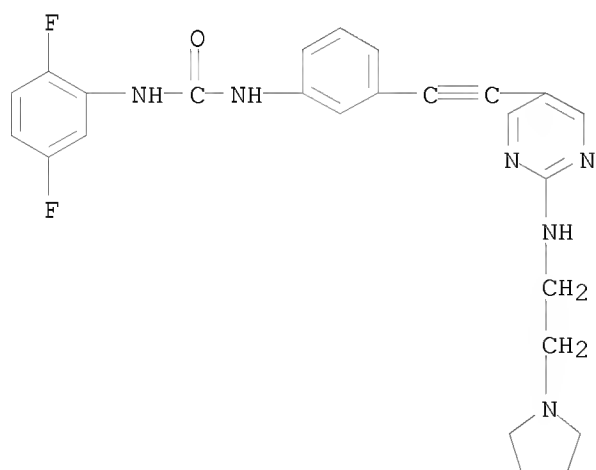
RN 857265-51-1 CAPLUS

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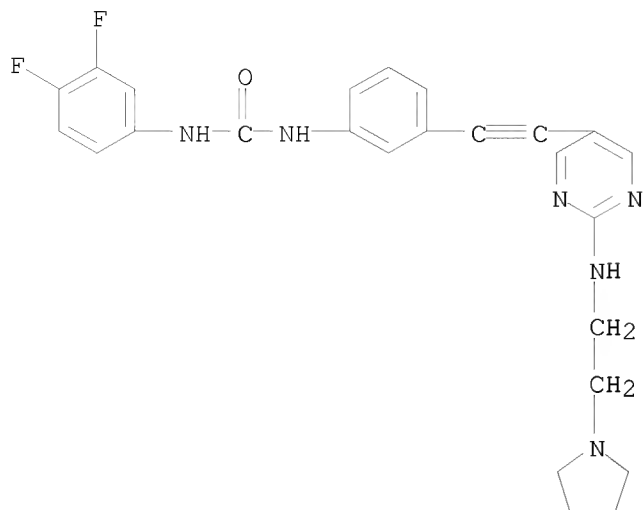
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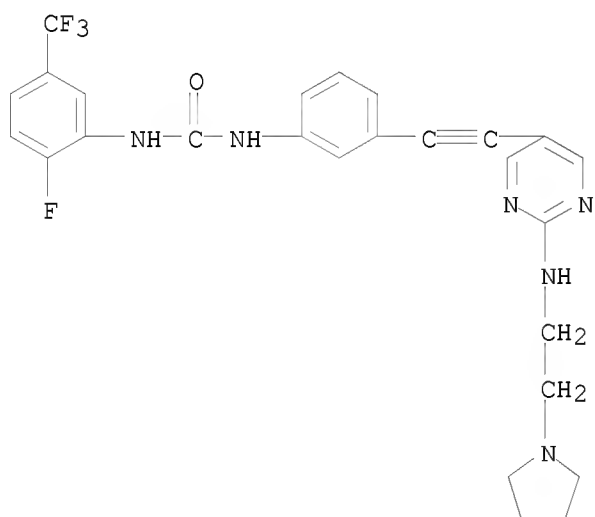
RN 857265-53-3 CAPLUS

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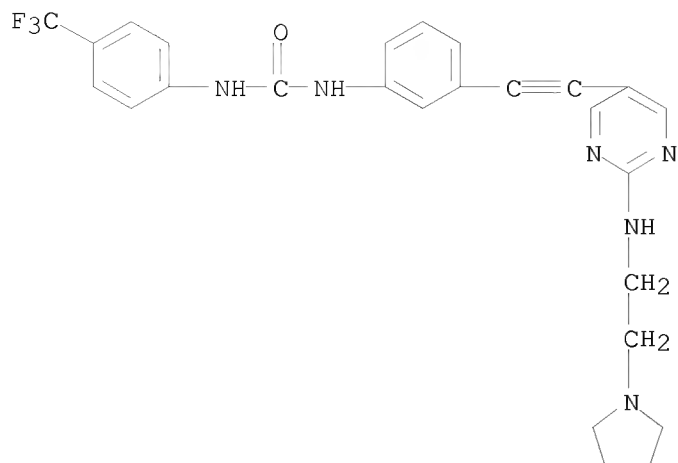
RN 857265-54-4 CAPLUS

CN Urea, N-[2-fluoro-5-(trifluoromethyl)phenyl]-N'-[3-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



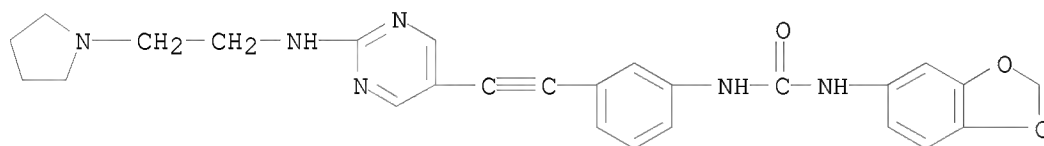
RN 857265-55-5 CAPLUS

CN Urea, N-[3-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[4-(trifluoromethyl)phenyl]- (CA INDEX NAME)



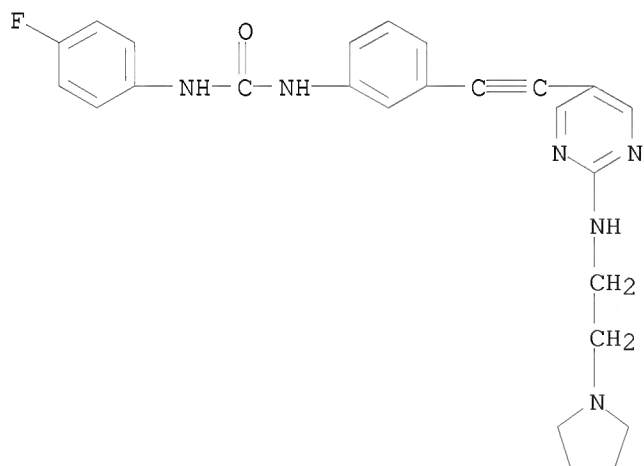
RN 857265-56-6 CAPLUS

CN Urea, N-1,3-benzodioxol-5-yl-N'-[3-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



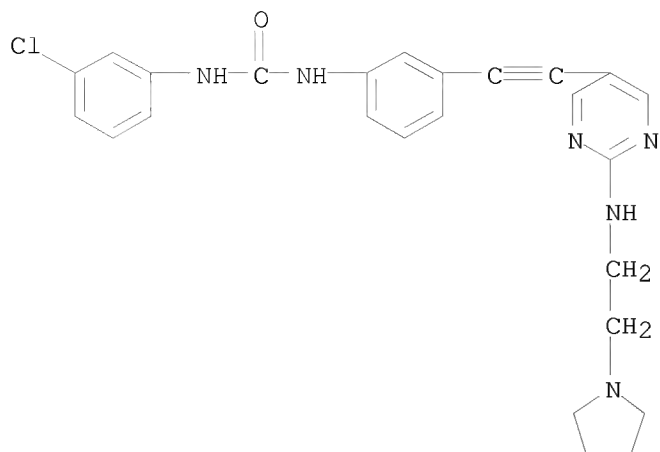
RN 857265-57-7 CAPLUS

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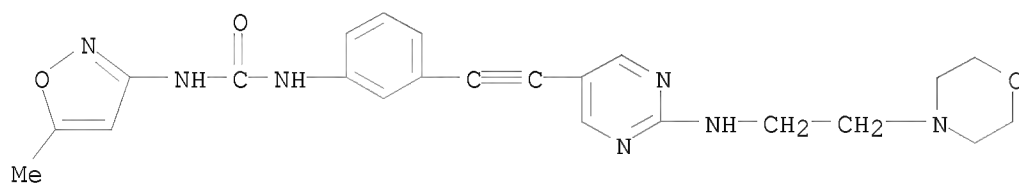


RN 857265-58-8 CAPLUS

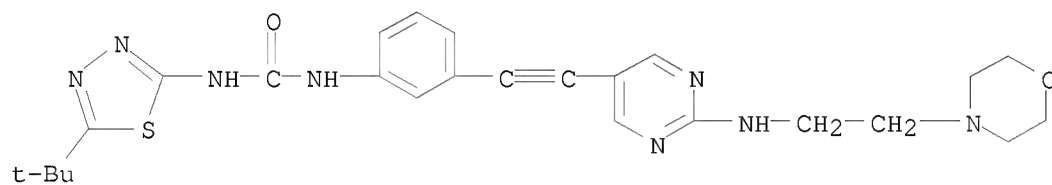
CN Urea, N-(3-chlorophenyl)-N'-[3-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



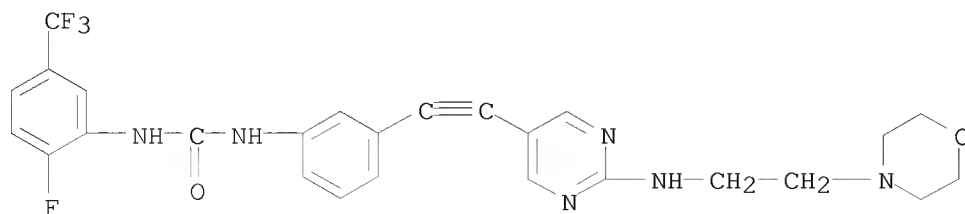
RN 857265-59-9 CAPLUS
 CN Urea, N-(5-methyl-3-isoxazolyl)-N'-[3-[2-[2-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857265-60-2 CAPLUS
 CN Urea, N-[5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl]-N'-[3-[2-[2-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

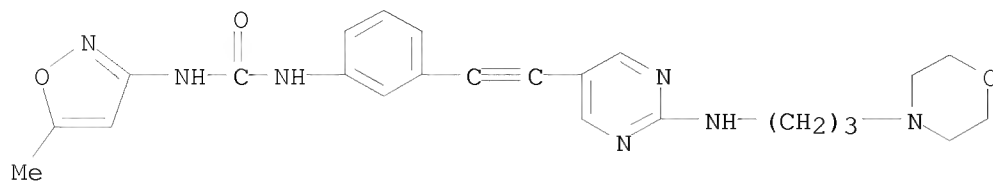


RN 857265-61-3 CAPLUS
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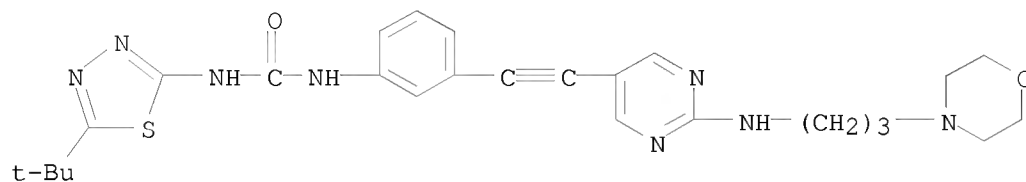
RN 857265-62-4 CAPLUS
 CN Urea, N-(5-methyl-3-isoxazolyl)-N'-[3-[2-[2-[[3-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-

morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



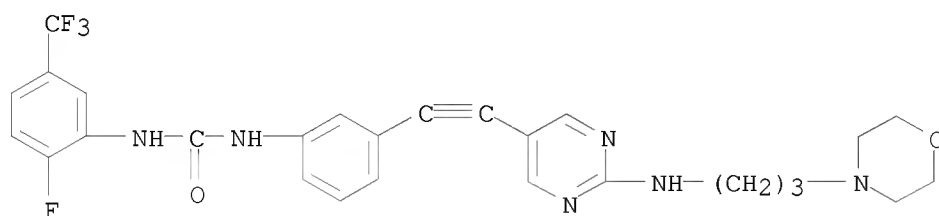
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CN Urea, N-[5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl]-N'-[3-[2-[2-[[3-(4-morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



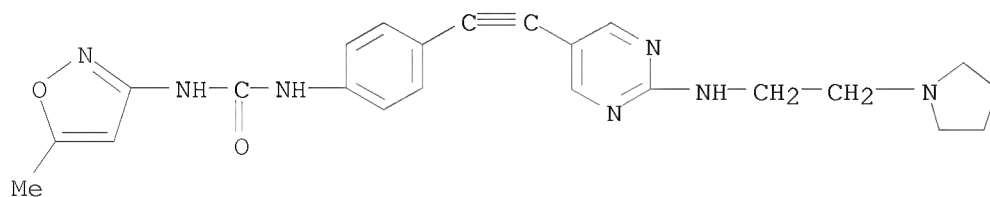
RN 857265-64-6 CAPLUS

CN Urea, N-[2-fluoro-5-(trifluoromethyl)phenyl]-N'-[3-[2-[2-[[3-(4-morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



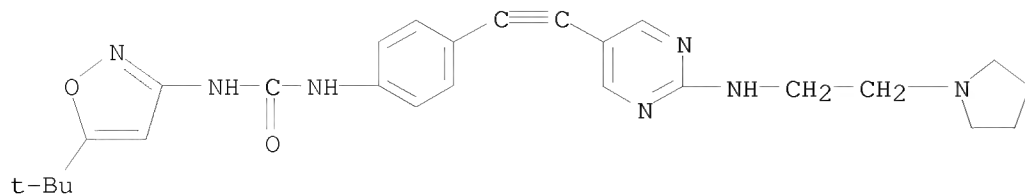
RN 857265-65-7 CAPLUS

CN Urea, N-(5-methyl-3-isoxazolyl)-N'-[4-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



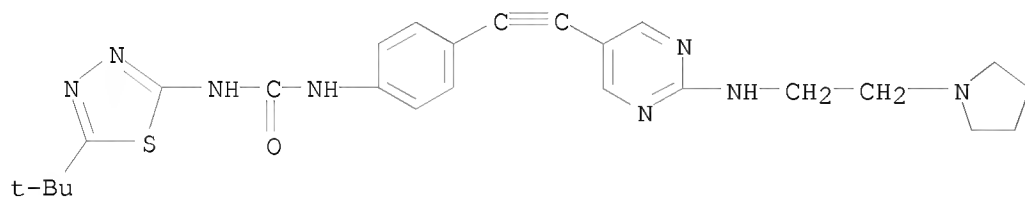
RN 857265-68-0 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[4-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



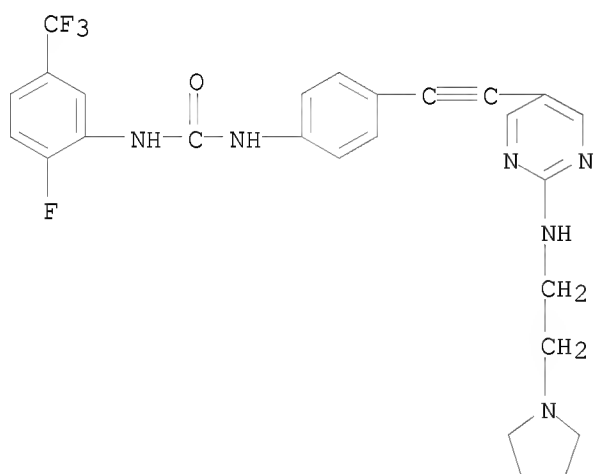
RN 857265-69-1 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl]-N'-[4-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857265-70-4 CAPLUS

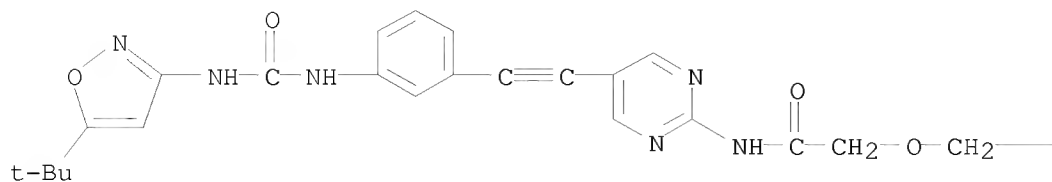
CN Urea, N-[2-fluoro-5-(trifluoromethyl)phenyl]-N'-[4-[2-[2-[[2-(1-pyrrolidinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857265-71-5 CAPLUS

CN Acetamide, N-[5-[2-[3-[[[5-(1,1-dimethylethyl)-3-isoxazolyl]amino]carbonyl]amino]phenyl]ethynyl]-2-(2-methoxyethoxy)- (CA INDEX NAME)

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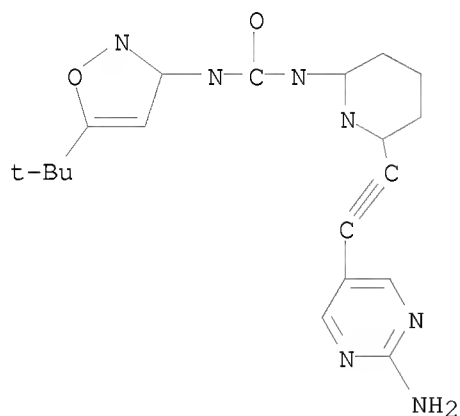


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RN 857265-72-6 CAPLUS

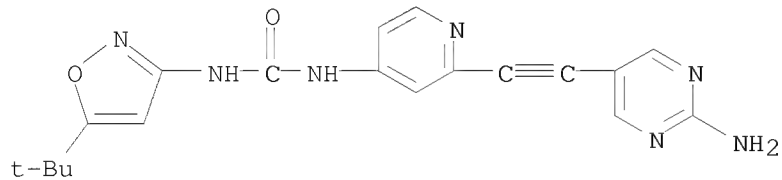
CN Urea, N-[6-[2-(2-amino-5-pyrimidinyl)ethynyl]-2-pyridinyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

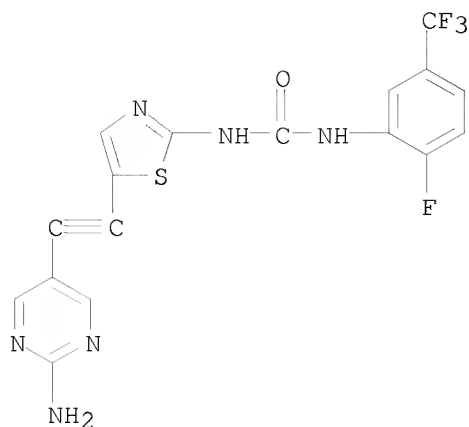
RN 857265-76-0 CAPLUS

CN Urea, N-[2-[2-(2-amino-5-pyrimidinyl)ethynyl]-4-pyridinyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



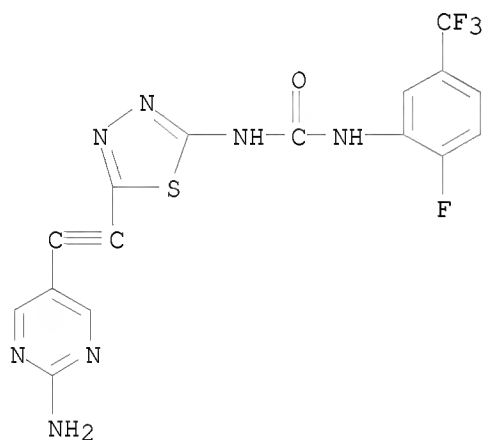
RN 857265-78-2 CAPLUS

CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-2-thiazolyl]-N'-[2-fluoro-5-(trifluoromethyl)phenyl]- (CA INDEX NAME)



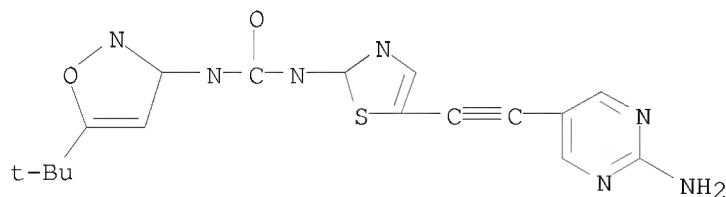
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CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-1,3,4-thiadiazol-2-yl]-N'-[2-fluoro-5-(trifluoromethyl)phenyl]- (CA INDEX NAME)



RN 857265-82-8 CAPLUS

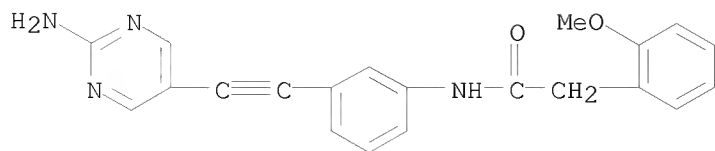
CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-2-thiazolyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



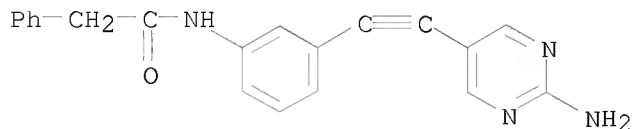
ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

RN 857265-84-0 CAPLUS

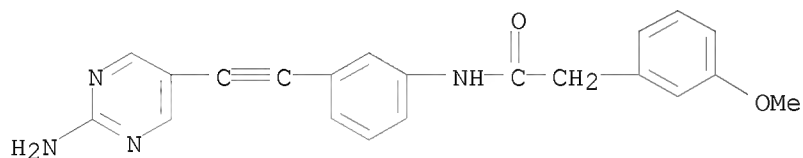
CN Benzeneacetamide, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-2-methoxy- (CA INDEX NAME)



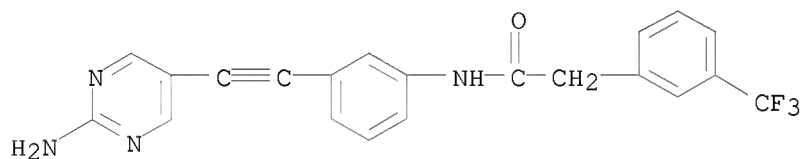
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 CN Benzeneacetamide, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]- (CA INDEX NAME)



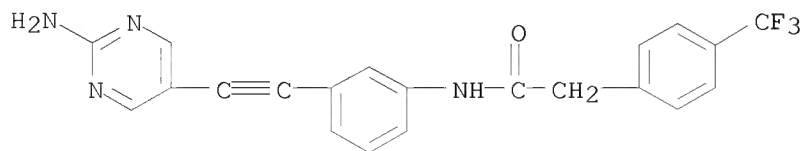
RN 857265-86-2 CAPLUS
 CN Benzeneacetamide, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-3-methoxy- (CA INDEX NAME)



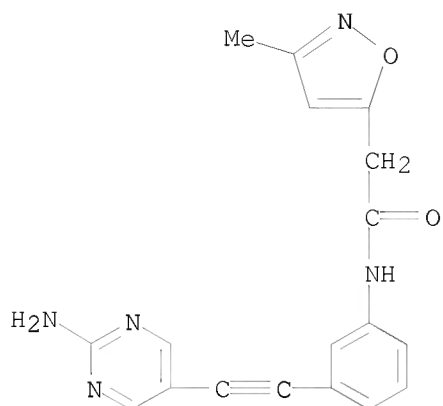
RN 857265-87-3 CAPLUS
 CN Benzeneacetamide, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-3-(trifluoromethyl)- (CA INDEX NAME)



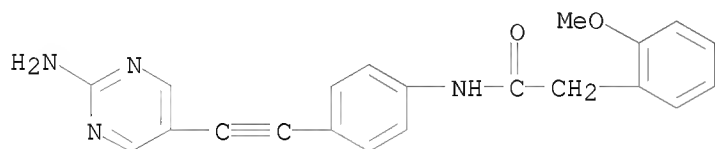
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 CN Benzeneacetamide, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-4-(trifluoromethyl)- (CA INDEX NAME)



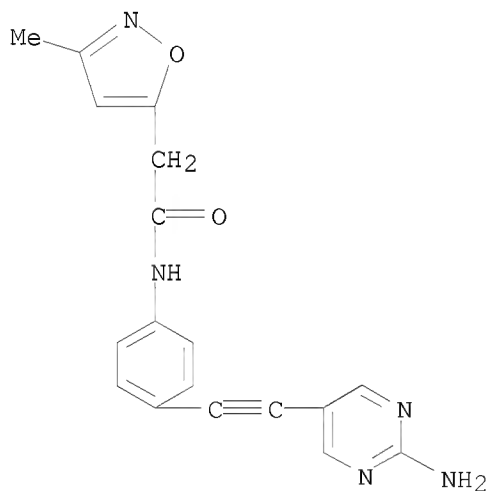
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 CN 5-Isoxazoleacetamide, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-3-methyl- (CA INDEX NAME)



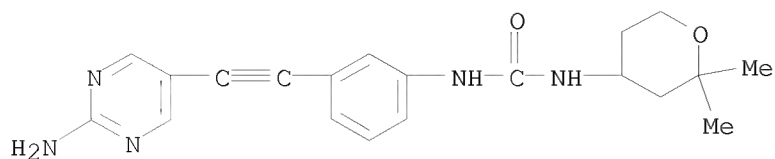
RN 857265-91-9 CAPLUS
 CN Benzeneacetamide, N-[4-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-2-methoxy-
 (CA INDEX NAME)



RN 857265-92-0 CAPLUS
 CN 5-Isoxazoleacetamide, N-[4-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-3-
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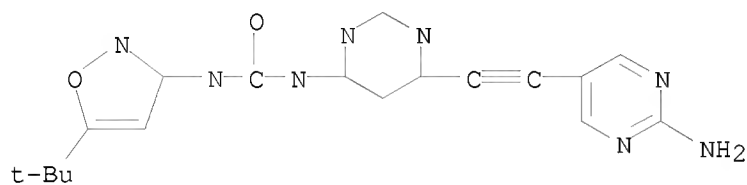


RN 857265-93-1 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(tetrahydro-2,2-
 dimethyl-2H-pyran-4-yl)- (CA INDEX NAME)



RN 857265-94-2 CAPLUS

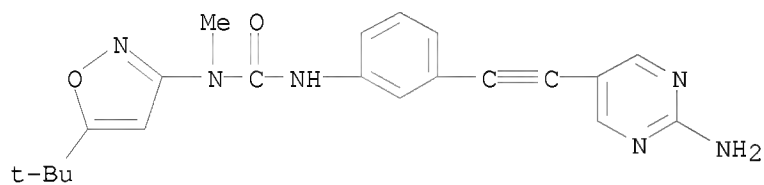
CN Urea, N-[6-[2-(2-amino-5-pyrimidinyl)ethynyl]-4-pyrimidinyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



ONE OR MORE TAUTOMERIC DOUBLE BONDS NOT DISPLAYED IN THE STRUCTURE

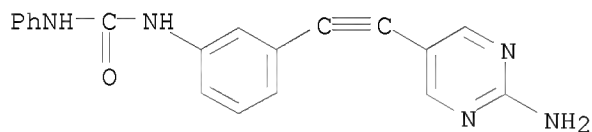
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CN Urea, N'-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N-methyl- (CA INDEX NAME)



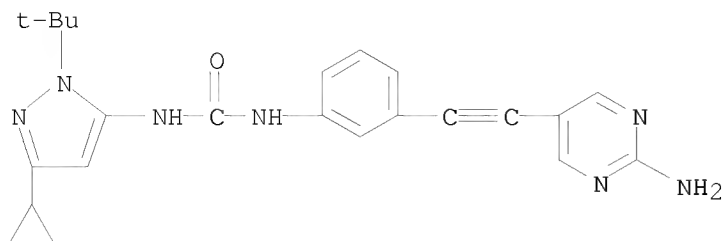
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CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-phenyl- (CA INDEX NAME)



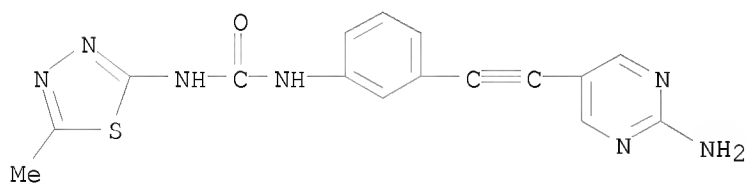
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CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[3-cyclopropyl-1-(1,1-dimethylethyl)-1H-pyrazol-5-yl]- (CA INDEX NAME)



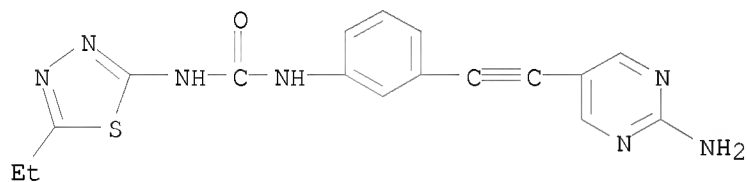
RN 857265-99-7 CAPLUS

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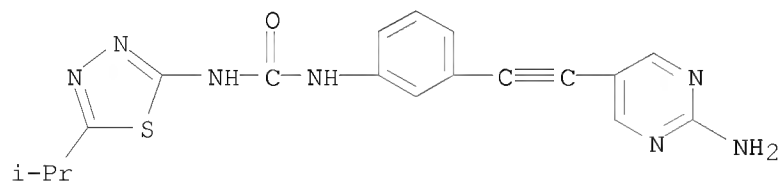
RN 857266-00-3 CAPLUS

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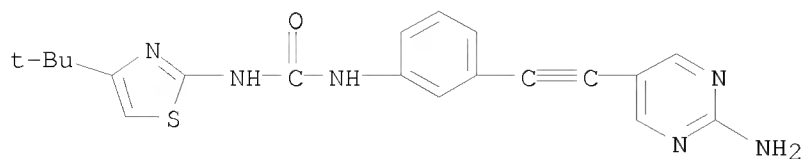
RN 857266-01-4 CAPLUS

CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(5-(1-methylethyl)-1,3,4-thiadiazol-2-yl)- (CA INDEX NAME)



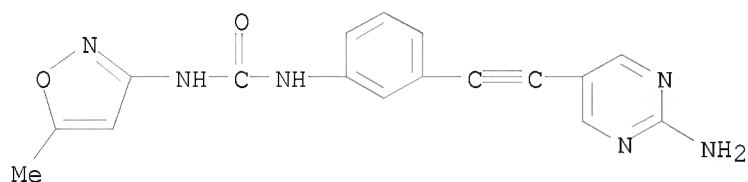
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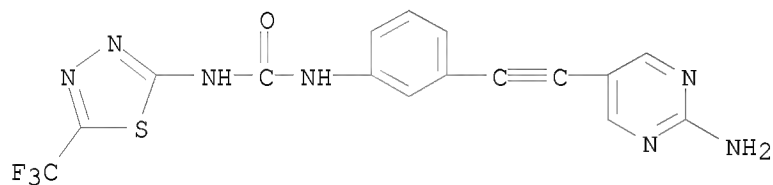
RN 857266-03-6 CAPLUS

CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(5-methyl-3-isoxazolyl)- (CA INDEX NAME)



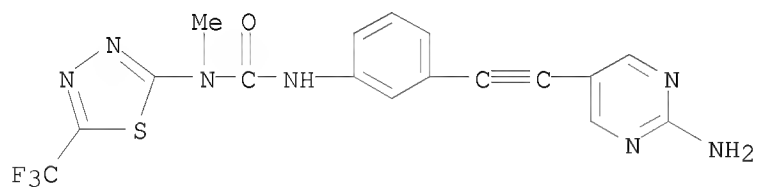
RN 857266-04-7 CAPLUS

CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[5-(trifluoromethyl)-1,3,4-thiadiazol-2-yl]- (CA INDEX NAME)



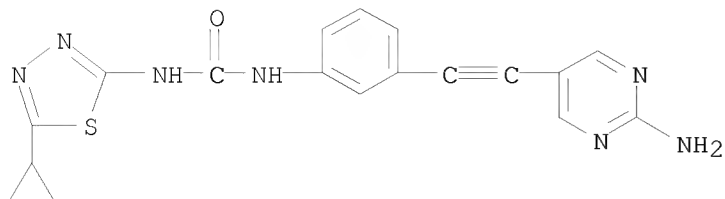
RN 857266-05-8 CAPLUS

CN Urea, N'-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N-methyl-N-[5-(trifluoromethyl)-1,3,4-thiadiazol-2-yl]- (CA INDEX NAME)



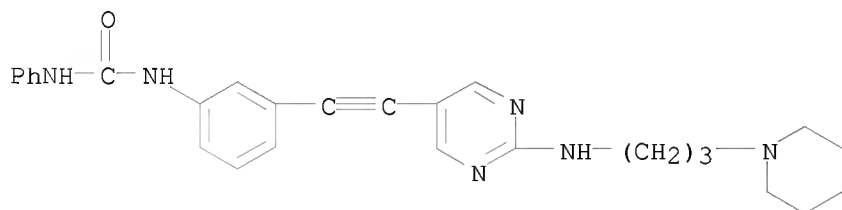
RN 857266-06-9 CAPLUS

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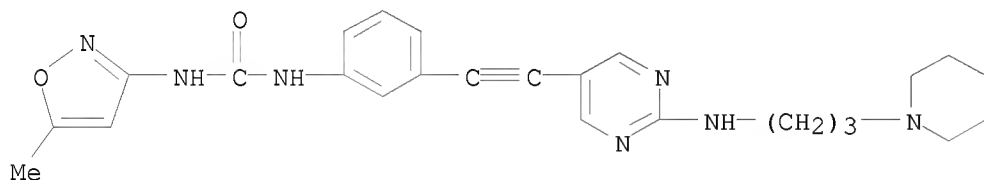
RN 857266-07-0 CAPLUS

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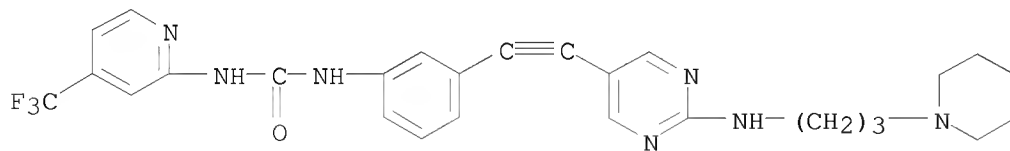
RN 857266-09-2 CAPLUS

CN Urea, N-(5-methyl-3-isoxazolyl)-N'-[3-[2-[2-[[3-(1-piperidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



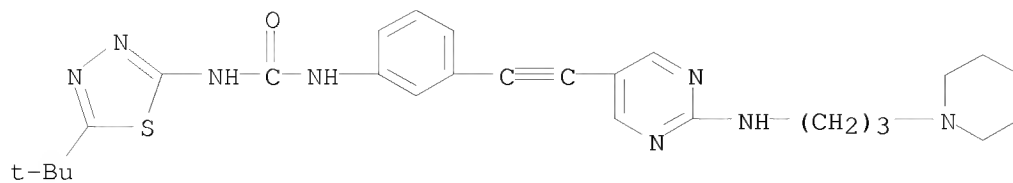
RN 857266-10-5 CAPLUS

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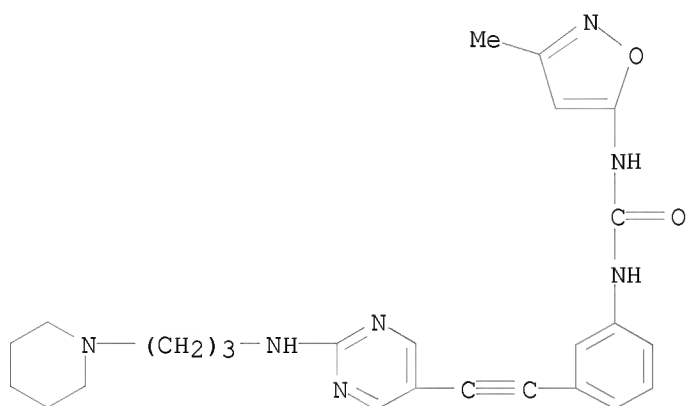
RN 857266-11-6 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-1,3,4-thiadiazol-2-yl]-N'-[3-[2-[2-[[3-(1-piperidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



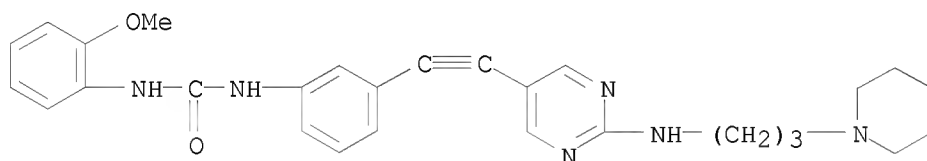
RN 857266-12-7 CAPLUS

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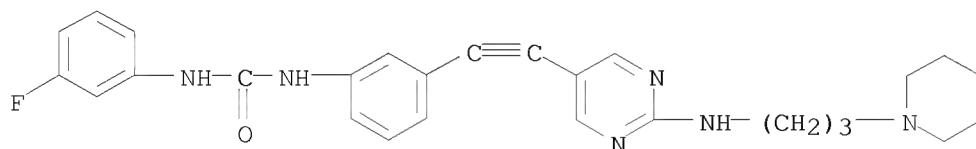
RN 857266-13-8 CAPLUS

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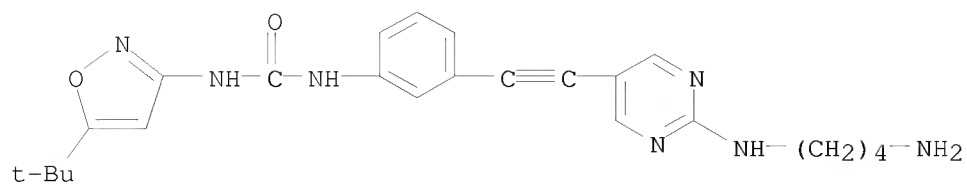
RN 857266-14-9 CAPLUS

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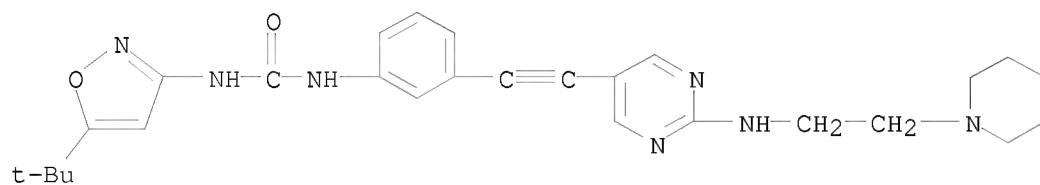
RN 857266-15-0 CAPLUS

CN Urea, N-[3-[2-[2-[[3-(1-piperidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



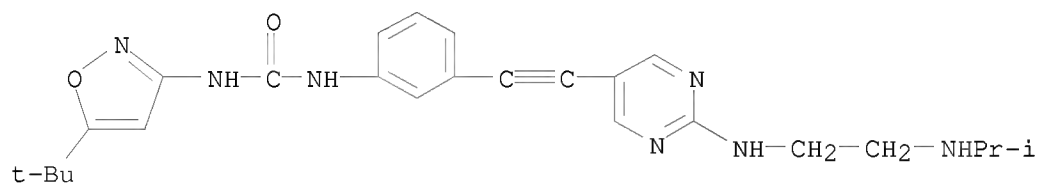
RN 857266-16-1 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[2-(1-piperidiny)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857266-17-2 CAPLUS

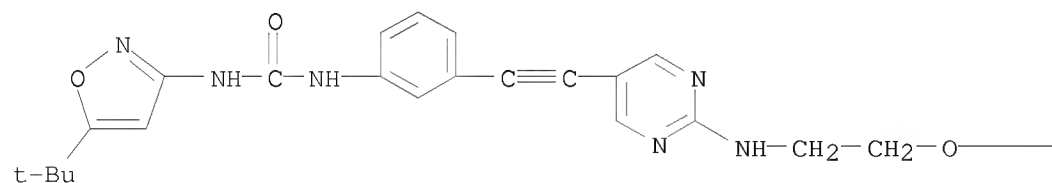
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[2-(1-methylethyl)amino]ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857266-18-3 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[2-(2-hydroxyethoxy)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

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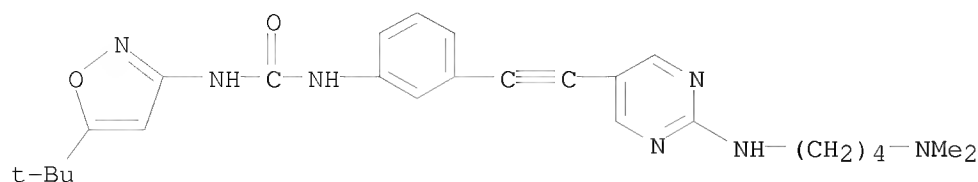


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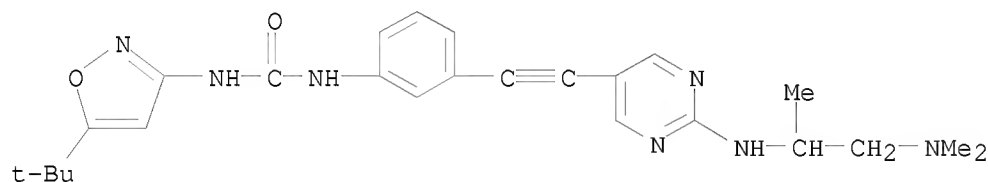
RN 857266-19-4 CAPLUS

CN Urea, N-[3-[2-[2-[[4-(dimethylamino)butyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



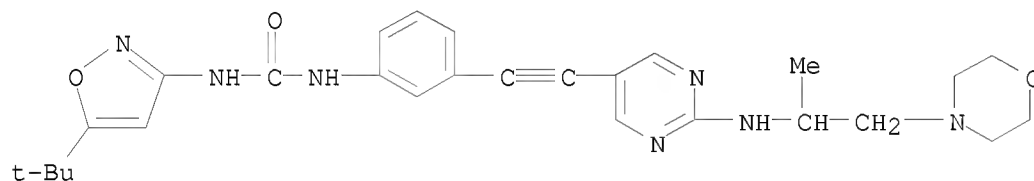
RN 857266-20-7 CAPLUS

CN Urea, N-[3-[2-[2-[[2-(dimethylamino)-1-methylethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)



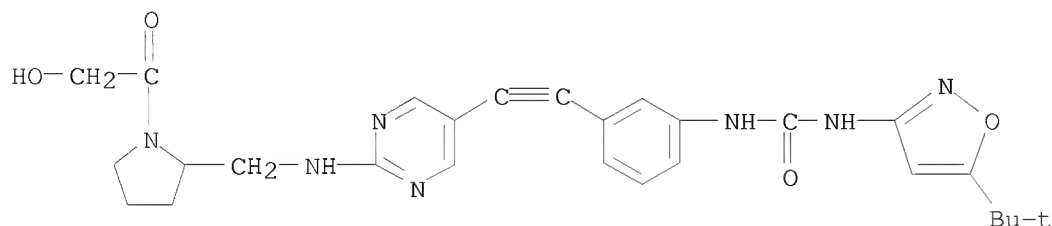
RN 857266-21-8 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[1-methyl-2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857266-22-9 CAPLUS

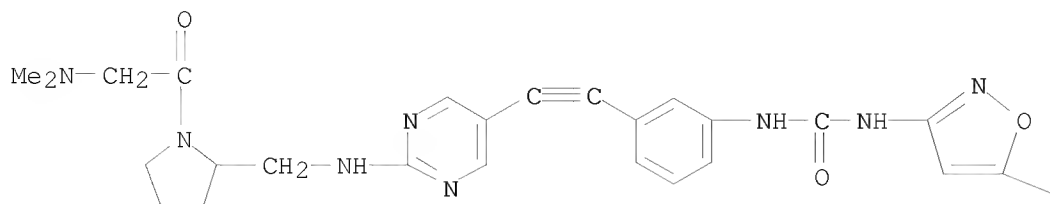
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[[1-(2-hydroxyacetyl)-2-pyrrolidinyl]methyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857266-25-2 CAPLUS

CN Urea, N-[3-[2-[2-[[[1-[2-(dimethylamino)acetyl]-2-pyrrolidinyl]methyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]- (CA INDEX NAME)

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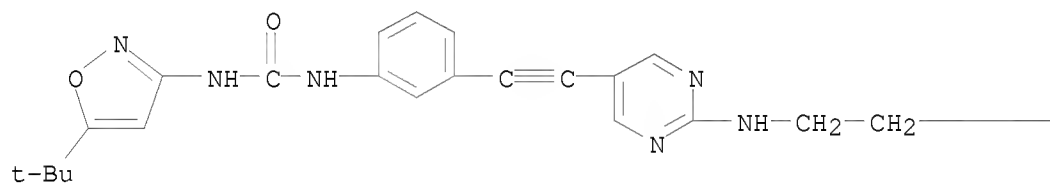


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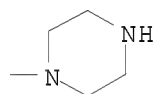
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RN 857266-26-3 CAPLUS
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[2-(1-piperazinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

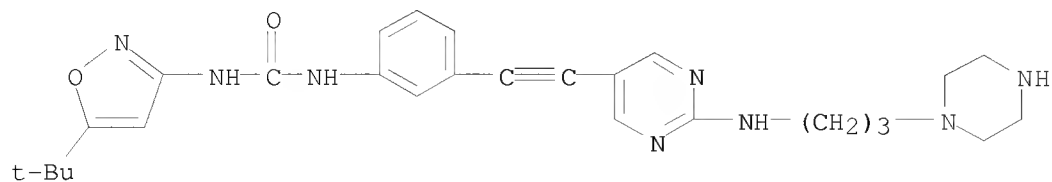
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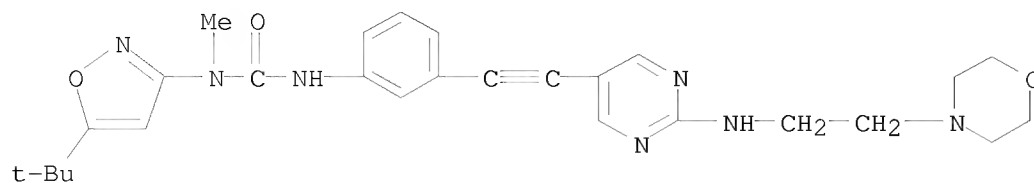


RN 857266-27-4 CAPLUS
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[[3-(1-piperazinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



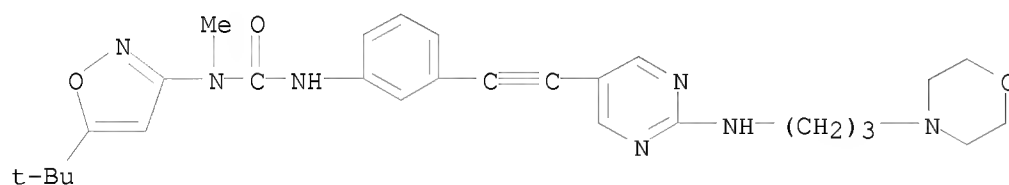
RN 857266-29-6 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N-methyl-N'-[3-[2-[2-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



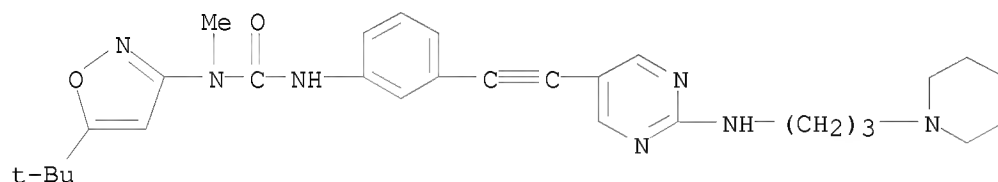
RN 857266-32-1 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N-methyl-N'-[3-[2-[2-[[3-(4-morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



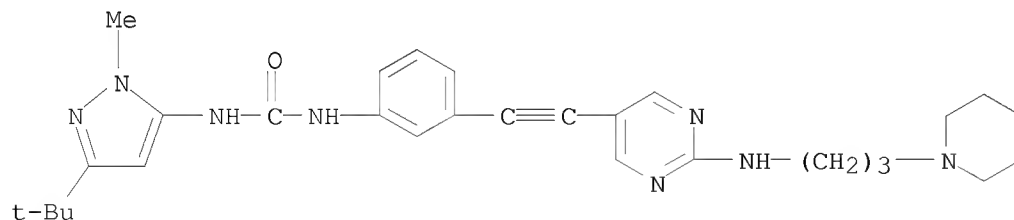
RN 857266-33-2 CAPLUS

CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N-methyl-N'-[3-[2-[2-[[3-(1-piperidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



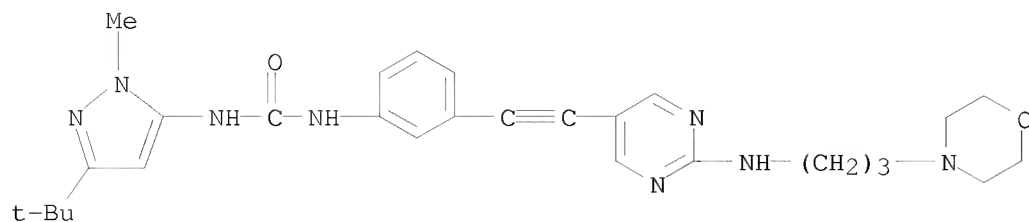
RN 857266-34-3 CAPLUS

CN Urea, N-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]-N'-[3-[2-[2-[[3-(1-piperidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



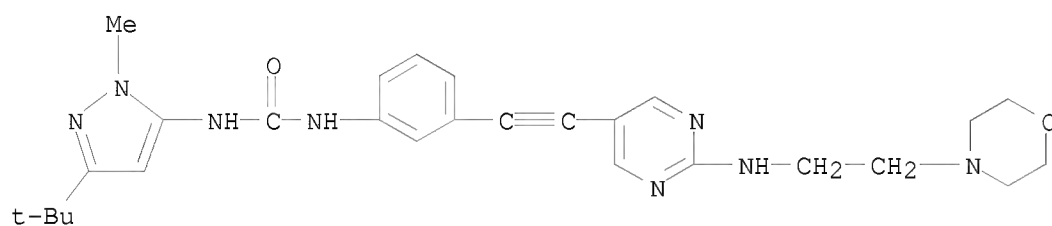
RN 857266-36-5 CAPLUS

CN Urea, N-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]-N'-[3-[2-[2-[[3-(4-morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



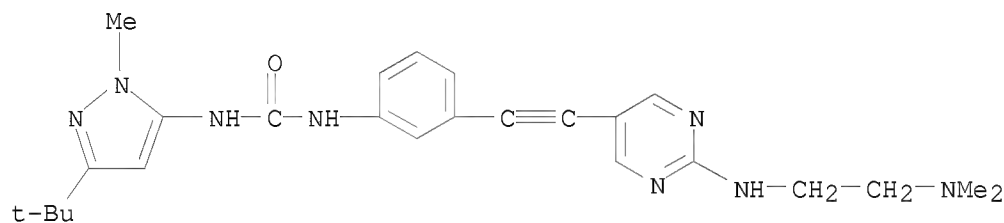
RN 857266-37-6 CAPLUS

CN Urea, N-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]-N'-[3-[2-[2-[[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



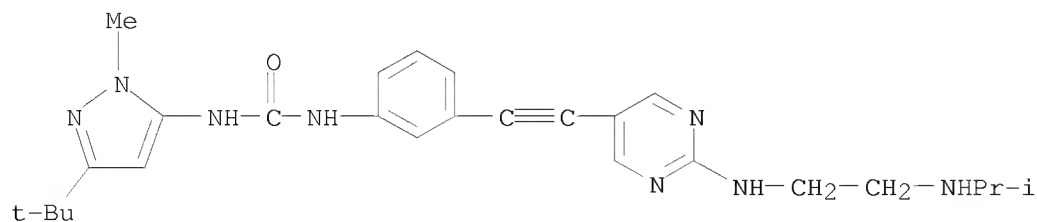
RN 857266-38-7 CAPLUS

CN Urea, N-[3-[2-[2-[[2-(dimethylamino)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]- (CA INDEX NAME)



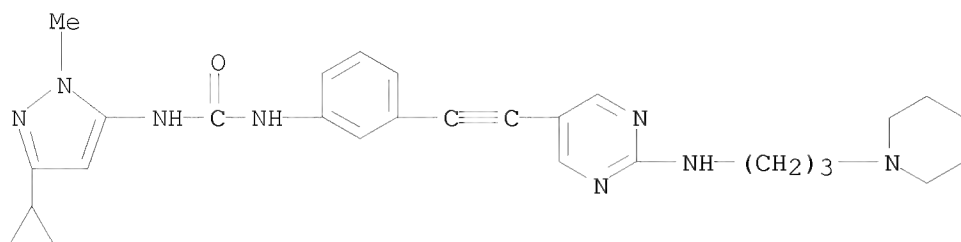
RN 857266-39-8 CAPLUS

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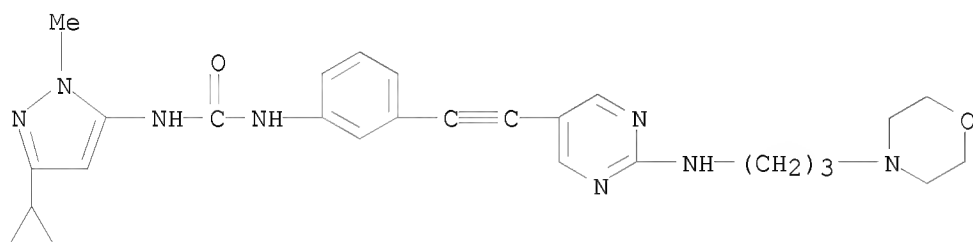


RN 857266-40-1 CAPLUS

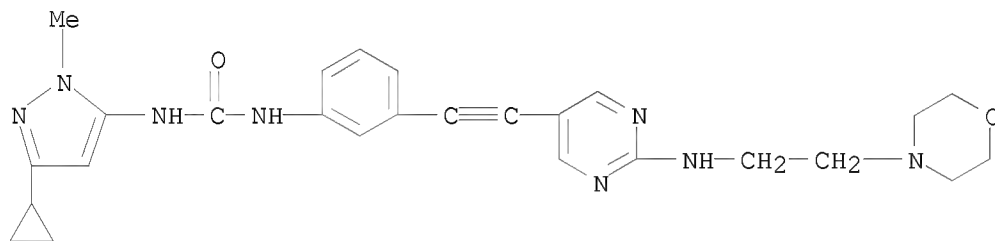
CN Urea, N-(3-cyclopropyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[2-[2-[[3-(1-piperidinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857266-42-3 CAPLUS
 CN Urea, N-(3-cyclopropyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[2-[2-[3-(4-morpholinyl)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)

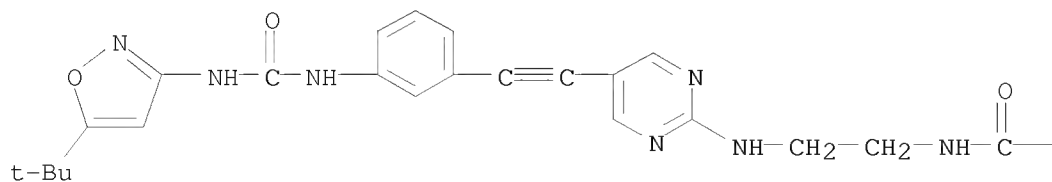


RN 857266-43-4 CAPLUS
 CN Urea, N-(3-cyclopropyl-1-methyl-1H-pyrazol-5-yl)-N'-[3-[2-[2-[2-(4-morpholinyl)ethyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857266-44-5 CAPLUS
 CN Acetamide, N-[2-[5-[2-[3-[[[[5-(1,1-dimethylethyl)-3-isoxazolyl]amino]carbonyl]amino]phenyl]ethynyl]-2-pyrimidinyl]amino]ethyl]-2-hydroxy- (CA INDEX NAME)

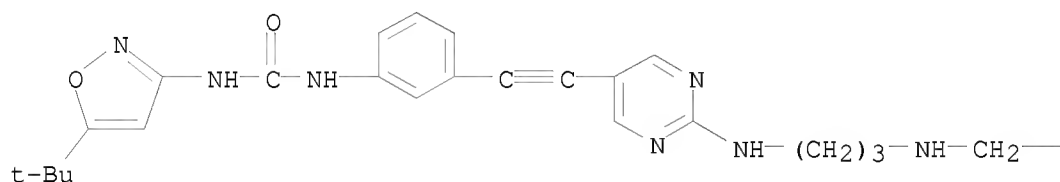
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RN 857266-45-6 CAPLUS

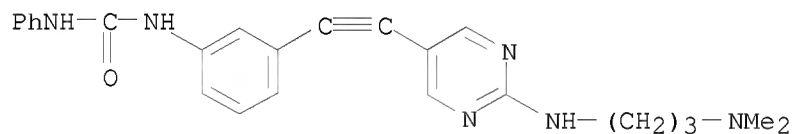
CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[3-[(2-hydroxyethyl)amino]propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



—CH₂—OH

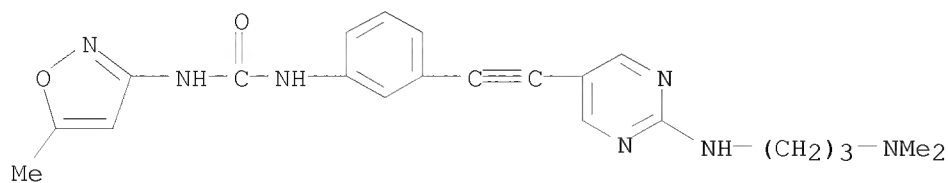
RN 857266-48-9 CAPLUS

CN Urea, N-[3-[2-[2-[3-(dimethylamino)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-phenyl- (CA INDEX NAME)



RN 857266-49-0 CAPLUS

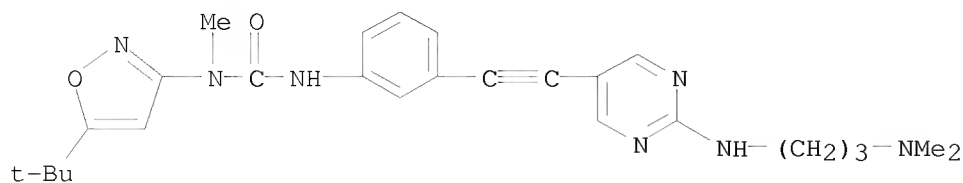
CN Urea, N-[3-[2-[2-[3-(dimethylamino)propyl]amino]-5-pyrimidinyl]ethynyl]phenyl]-N'-(5-methyl-3-isoxazolyl)- (CA INDEX NAME)



RN 857266-50-3 CAPLUS

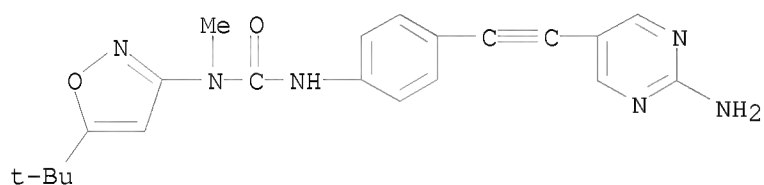
CN Urea, N'-[3-[2-[2-[3-(dimethylamino)propyl]amino]-5-

pyrimidinyl]ethynyl]phenyl]-N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N-methyl- (CA INDEX NAME)



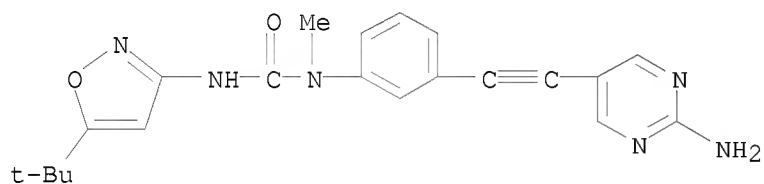
RN 857266-51-4 CAPLUS

CN Urea, N'-[4-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N-methyl- (CA INDEX NAME)



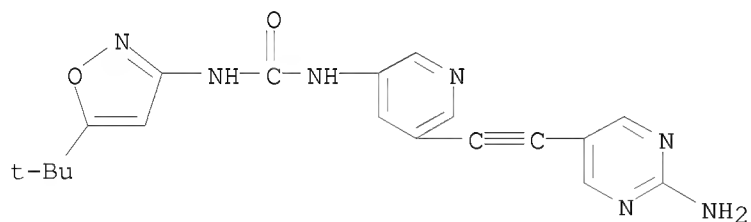
RN 857266-53-6 CAPLUS

CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N-methyl- (CA INDEX NAME)



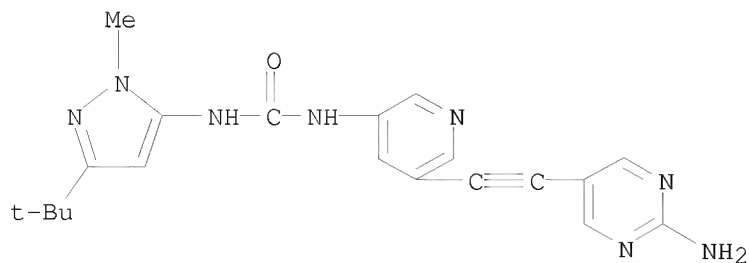
RN 857266-57-0 CAPLUS

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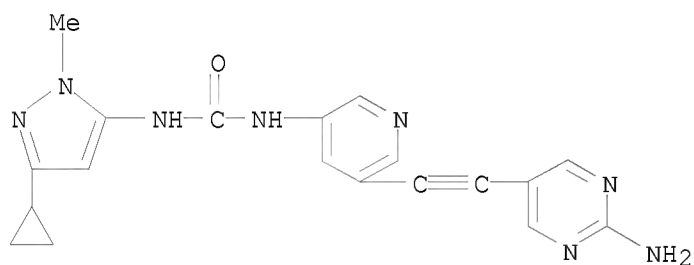
RN 857266-61-6 CAPLUS

CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-3-pyridinyl]-N'-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]- (CA INDEX NAME)



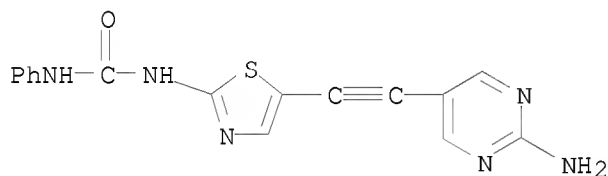
RN 857266-63-8 CAPLUS

CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-3-pyridinyl]-N'-(3-cyclopropyl-1-methyl-1H-pyrazol-5-yl)- (CA INDEX NAME)



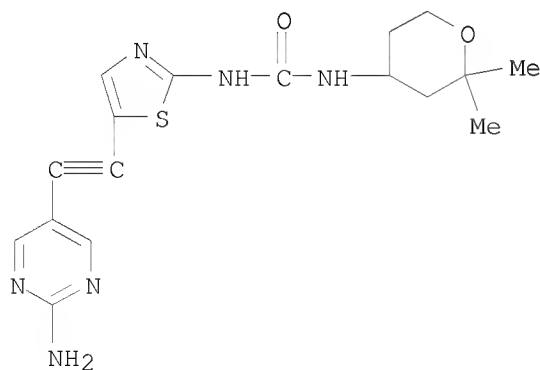
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CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-2-thiazolyl]-N'-phenyl- (CA INDEX NAME)



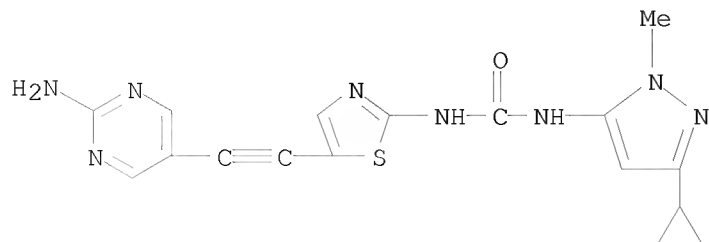
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CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-2-thiazolyl]-N'-(tetrahydro-2,2-dimethyl-2H-pyran-4-yl)- (CA INDEX NAME)



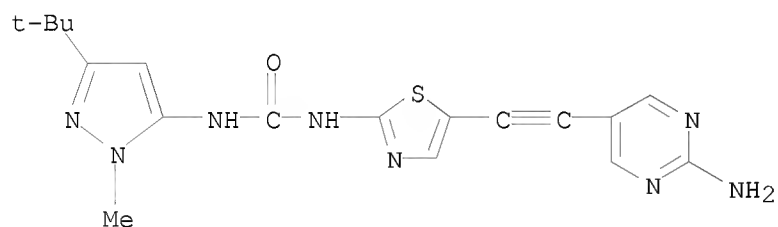
RN 857266-67-2 CAPLUS

CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-2-thiazolyl]-N'-(3-cyclopropyl-1-methyl-1H-pyrazol-5-yl)- (CA INDEX NAME)



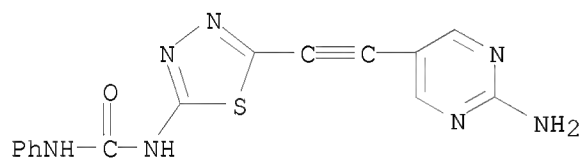
RN 857266-70-7 CAPLUS

CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-2-thiazolyl]-N'-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]- (CA INDEX NAME)



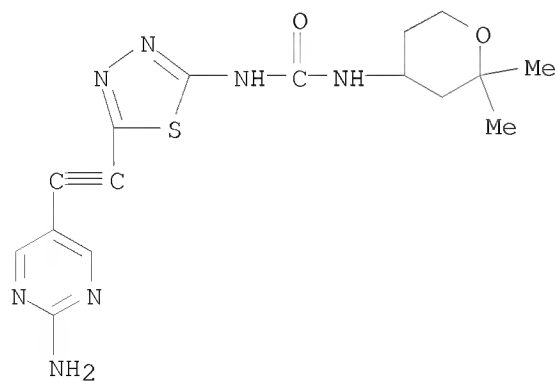
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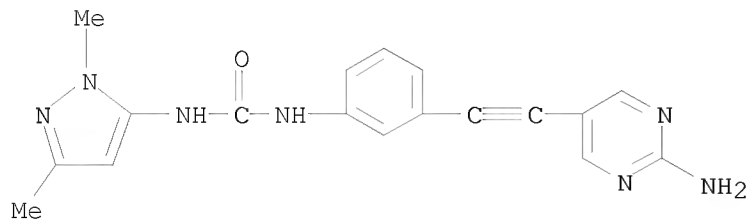


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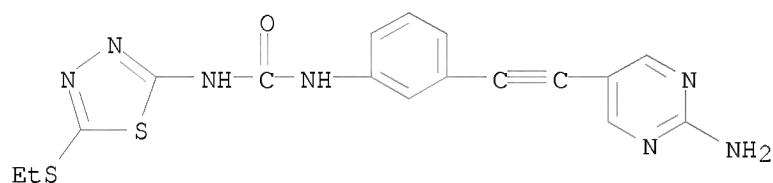
CN Urea, N-[5-[2-(2-amino-5-pyrimidinyl)ethynyl]-1,3,4-thiadiazol-2-yl]-N'-(tetrahydro-2,2-dimethyl-2H-pyran-4-yl)- (CA INDEX NAME)



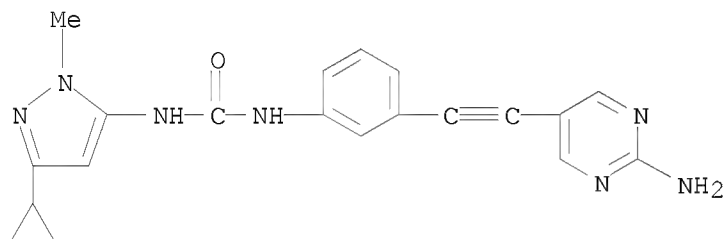
RN 857266-82-1 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(1,3-dimethyl-1H-pyrazol-5-yl)- (CA INDEX NAME)



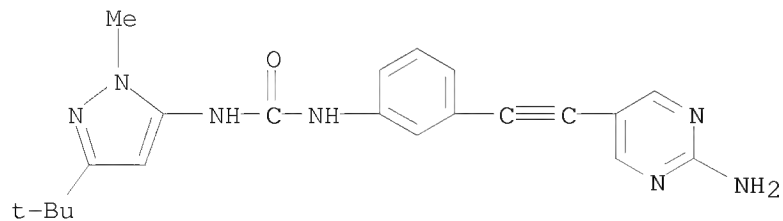
RN 857266-84-3 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[5-(ethylthio)-1,3,4-thiadiazol-2-yl]- (CA INDEX NAME)



RN 857266-86-5 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(3-cyclopropyl-1-methyl-1H-pyrazol-5-yl)- (CA INDEX NAME)

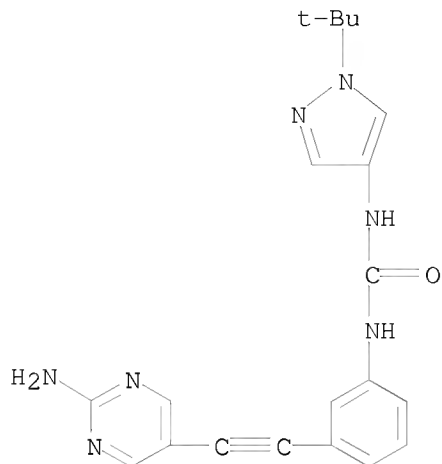


RN 857266-88-7 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]- (CA INDEX NAME)



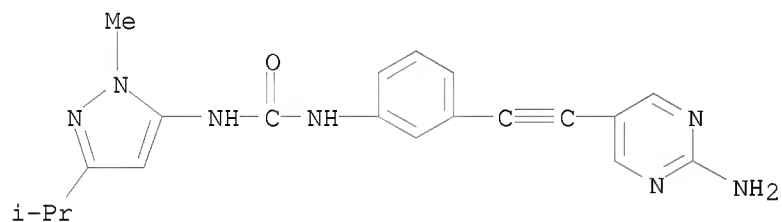
RN 857266-90-1 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[1-(1,1-

dimethylethyl)-1H-pyrazol-4-yl]- (CA INDEX NAME)



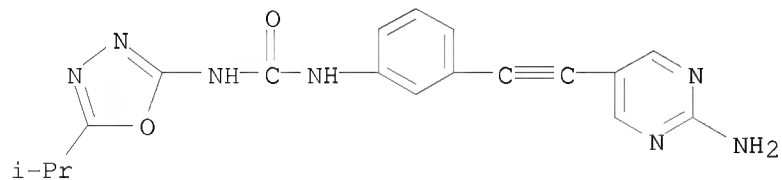
RN 857266-93-4 CAPLUS

CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[1-methyl-3-(1-methylethyl)-1H-pyrazol-5-yl]- (CA INDEX NAME)



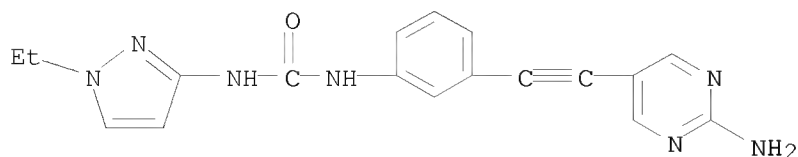
RN 857266-96-7 CAPLUS

CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[5-(1-methylethyl)-1,3,4-oxadiazol-2-yl]- (CA INDEX NAME)

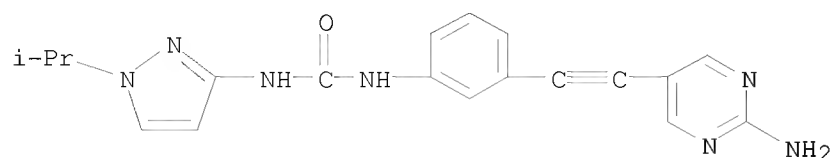


RN 857266-99-0 CAPLUS

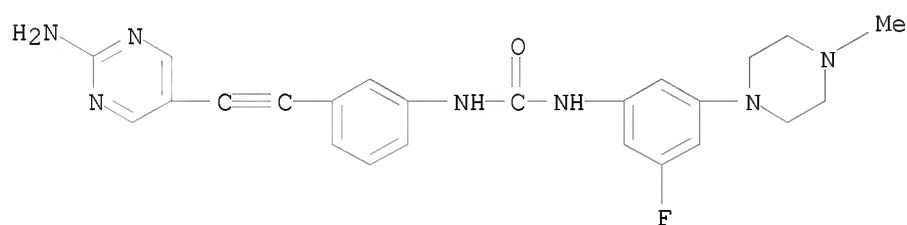
CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-(1-ethyl-1H-pyrazol-3-yl)- (CA INDEX NAME)



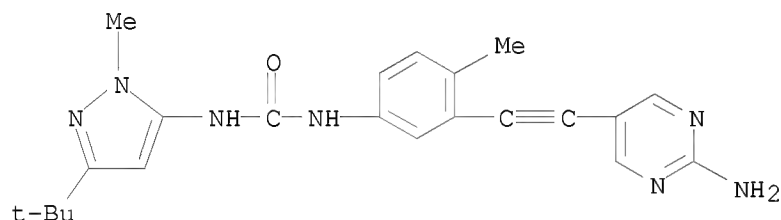
RN 857267-02-8 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[1-(1-methylethyl)-1H-pyrazol-3-yl]- (CA INDEX NAME)



RN 857267-06-2 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]phenyl]-N'-[3-fluoro-5-(4-methyl-1-piperazinyl)phenyl]- (CA INDEX NAME)

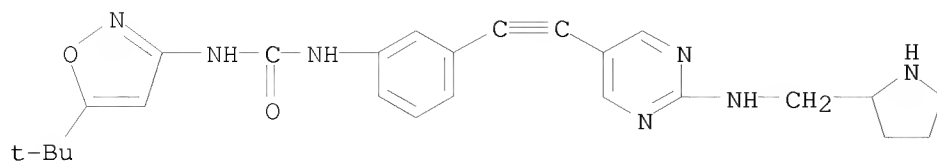


RN 857267-09-5 CAPLUS
 CN Urea, N-[3-[2-(2-amino-5-pyrimidinyl)ethynyl]-4-methylphenyl]-N'-[3-(1,1-dimethylethyl)-1-methyl-1H-pyrazol-5-yl]- (CA INDEX NAME)

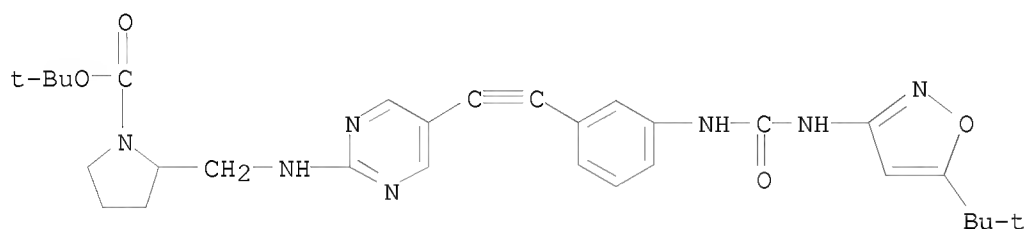


IT 857266-23-0P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[[pyrrolidin-2-yl)methyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea
 857266-24-1P, N-(5-tert-Butylisoxazol-3-yl)-N'-[3-[[2-[[[1-(tert-butoxycarbonyl)pyrrolidin-2-yl)methyl]amino]pyrimidin-5-yl]ethynyl]phenyl]urea 857266-52-5P, Phenyl [4-[(2-aminopyrimidin-5-yl)ethynyl]phenyl]carbamate
 RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
 (preparation of pyrimidine derivs. as inhibitors of Tie2 receptor tyrosine kinases)

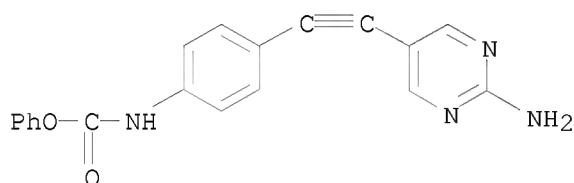
RN 857266-23-0 CAPLUS
 CN Urea, N-[5-(1,1-dimethylethyl)-3-isoxazolyl]-N'-[3-[2-[2-[(2-pyrrolidinylmethyl)amino]-5-pyrimidinyl]ethynyl]phenyl]- (CA INDEX NAME)



RN 857266-24-1 CAPLUS
 CN 1-Pyrrolidinecarboxylic acid, 2-[[[5-[2-[3-[[[5-(1,1-dimethylethyl)-3-isoxazolyl]amino]carbonyl]amino]phenyl]ethynyl]-2-pyrimidinyl]amino]methyl]-, 1,1-dimethylethyl ester (CA INDEX NAME)



RN 857266-52-5 CAPLUS
 CN Carbamic acid, [4-[(2-amino-5-pyrimidinyl)ethynyl]phenyl]-, phenyl ester (9CI) (CA INDEX NAME)



REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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LOGOFF? (Y)/N/HOLD:y

COST IN U.S. DOLLARS

SINCE FILE TOTAL

ENTRY SESSION

FULL ESTIMATED COST

16.83 196.80

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE TOTAL

ENTRY SESSION

CA SUBSCRIBER PRICE

-2.40 -2.40

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